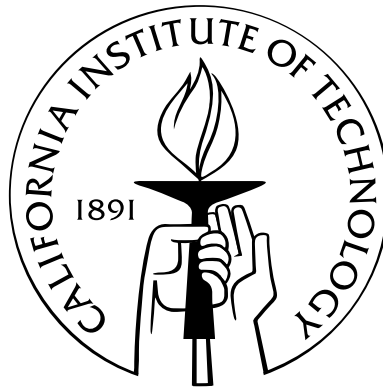


SUPERSTRING HOLOGRAPHY AND INTEGRABILITY IN $AdS_5 \times S^5$

Dissertation by

Ian Swanson

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For my parents, Frank and Sheila, and my brother, Cory.

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Abstract

The AdS/CFT correspondence provides a rich testing ground for many important topics in theoretical physics. The earliest and most striking example of the correspondence is the conjectured duality between the energy spectrum of type IIB superstring theory on $AdS_5 \times S^5$ and the operator anomalous dimensions of $\mathcal{N} = 4$ supersymmetric Yang-Mills theory in four dimensions. While there is a substantial amount of evidence in support of this conjecture, direct tests have been elusive. The difficulty of quantizing superstring theory in a curved Ramond-Ramond background is compounded by the problem of computing anomalous dimensions for non-BPS operators in the strongly coupled regime of the gauge theory. The former problem can be circumvented to some extent by taking a Penrose limit of $AdS_5 \times S^5$, reducing the background to that of a pp-wave (where the string theory is soluble). A corresponding limit of the gauge theory was discovered by Berenstein, Maldacena and Nastase, who obtained successful agreement between a class of operator dimensions in this limit and corresponding string energies in the Penrose limit. In this dissertation we present a body of work based largely on the introduction of worldsheet interaction corrections to the free pp-wave string theory by lifting the Penrose limit of $AdS_5 \times S^5$. This provides a new class of rigorous tests of AdS/CFT that probe a truly quantum realm of the string theory. By studying the correspondence in greater detail, we stand to learn not only about how the duality is realized on a more microscopic level, but how Yang-Mills theories behave at strong coupling. The methods presented here will hopefully contribute to the realization of these important goals.

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Introduction and overview

Since conservation laws arise from symmetries of the Lagrangian [1], an efficient way to characterize physical systems is to describe the mathematical symmetries under which they are invariant. From a certain perspective, the symmetries themselves may be viewed as paramount: a complete description of fundamental physics will likely be founded on an account of which symmetries are allowed by nature, under what circumstances these symmetries are realized and how and when these symmetries are broken. At the energies probed by current experiments, nature is described at the microscopic level by a quantum field theory with certain gauge symmetries. This framework is remarkably successful at describing particle spectra and interactions, but there are many convincing indications that this picture breaks down near the Planck scale, where gravitational effects become important.

To incorporate gravity in a way that is consistent at the quantum level, one must make a dramatic departure from the point-particle quantum field theory upon which the Standard Model is based. Only by replacing the fundamental point-particle constituents of the theory with one-dimensional extended objects (strings) is one afforded the freedom necessary to accommodate gravity [2, 3]. The physical theory of these objects, or string theory, is not only able to provide a consistent theory of quantum gravity, but also has a rich enough structure to give rise to the types of gauge symmetries observed in nature (and is free of quantum anomalies) [2–7]. One fascinating aspect of string theory, however, is that quantum consistency demands that the theory occupies ten spacetime dimensions (M-theory is eleven dimensional). Since we observe only four spacetime dimensions in the universe, theorists are charged

with the task of understanding the role of the six ‘extra’ spatial dimensions that are predicted by string theory. At first glance, the idea that six spatial dimensions exist in the universe but are somehow hidden seems fanciful. Stated concisely, a strong hope among theorists is that the extra dimensions in string theory will provide a mechanism through which the gauge symmetries in nature are realized naturally.

In the course of trying to describe the known symmetries of the vacuum, the study of string theory has led to the discovery of a dramatically new class of fundamental symmetries known as dualities. These symmetries stand apart from more traditional examples in that they connect physical theories that, at least superficially, appear to be entirely distinct in their formulation. This notion of duality, or the underlying equivalence of two seemingly disparate physical systems, has emerged as a powerful tool in recent decades. The usefulness of duality derives in part from the fact that dual descriptions are typically complementary, insofar as information that is inaccessible in one physical theory may often be extracted from a straightforward calculation in the theory’s dual description. This is often realized in the form of a strong/weak duality, whereby a small parameter useful for perturbation theory on one side is mapped to a large parameter on the other. Information provided by a perturbative expansion in one theory therefore equates to knowledge about nonperturbative physics in the dual theory (and vice versa).

In this work we will primarily be concerned with dualities that arise holographically, meaning that information (or degrees of freedom) existing in one theory with a given number of spacetime dimensions can be encoded in some dual theory with fewer spatial dimensions. This is of course analogous to an actual hologram, wherein information about the shape of an object in three spatial dimensions can be encoded on a two-dimensional film: in addition to recording the location in two dimensions of laser light incident on its surface, a hologram records the polarization of this light as it is reflected off of the object. A major theme in holographic dualities is that the importance of the spatial dimensions in which a theory is defined is often secondary to a proper accounting of the degrees of freedom accessible to the theory. This leads

us to how holography was initially recognized as an important concept in theoretical physics: the black-hole entropy problem.

0.1 The holographic entropy bound

As described above, the degrees of freedom in the universe appear to be described by quantum fields living in a four-dimensional spacetime, at least down to the scales accessible to current accelerator experiments. The belief among theorists is that this description holds all the way down to the Planck scale, l_{Planck} . The implication is that, with l_{Planck} serving as an ultraviolet cutoff, the degrees of freedom available to the vacuum can be roughly described by a three-dimensional lattice theory with internal lattice spacing equal to l_{Planck} . With one binary degree of freedom per Planck volume, the maximum entropy of a system enclosed in a volume V should scale in direct proportion to V [8–10].

The limitations of this simple picture can be seen by considering a thermodynamic system in which gravitational effects are important: namely, a black hole. The entropy of an isolated black hole is given by the Bekenstein-Hawking formula [11, 12]:

$$S_{\text{BH}} = \frac{A}{4G} . \tag{0.1.1}$$

The most striking aspect of this formula is that S_{BH} scales linearly with the area A of the event horizon. A simple thought experiment, following Bekenstein [12–14], leads to an interesting problem. Imagine some volume V of space that contains a thermodynamic system with entropy $S > S_{\text{BH}}$. If the entropy of the system is bounded by its volume, then this is a reasonable proposal. The mass of the system must be no greater than the mass of a black hole whose horizon is the boundary of V , otherwise the system would be larger than V . Now, if a thin shell of mass collapses into the system and forms a black hole whose horizon is precisely defined by V , the entropy of the new system is given by the Bekenstein-Hawking formula: this process

violates the second law of thermodynamics.

A striking solution to this problem, proposed by 't Hooft [15], is that nature obeys a *holographic entropy bound*, which states that the degrees of freedom available to a physical system occupying a volume V can be mapped to some physical theory defined to exist strictly on the boundary ∂V (see also [8, 9, 16, 17]). The maximum entropy of a system is thus limited by the number of degrees of freedom that can be mapped from the interior of the system to its boundary. The most striking aspect of this claim is that, while both theories must give rise to equivalent physical predictions, the ‘dual’ theory defined on the boundary necessarily exists in a fewer number of spatial dimensions than the original theory living in the bulk.

0.2 Holography and string theory

The holographic principal is deeply enmeshed in the intricate relationship between string theory and point-particle gauge theory. As a toy example, consider the analogy between the classical statistical mechanics of a D dimensional system and the quantum dynamics of a $D - 1$ dimensional system. (This analogy was alluded to extensively by Polyakov in [18].) The statement for $D = 1$ is that the quantum transition amplitude for a point particle over some time interval T can be interpreted as the classical partition function of a string whose length is determined by T . Although not strictly holographic, this example captures several themes that are ubiquitous in gauge/string-theory dualities.

We should first take note of the types of gauge theories that will be of interest to us. The theory of the strong nuclear force, or quantum chromodynamics (QCD), is an $SU(3)$ gauge theory: it is a non-Abelian Yang-Mills theory with three colors ($N_c = 3$). QCD is known to be asymptotically free, meaning that the theory is free at high energies. At very low energies one enters a regime where perturbation theory is no longer useful, and with no further advancements (such as a dual string formulation) the only hope is that lattice computations will one day be able to probe

these regions of the theory in detail. In 1974 't Hooft suggested that a more general $SU(N_c)$ Yang-Mills theory would simplify when the rank of the gauge group (or the number of colors) N_c becomes large [19]. Such a simplification is intriguing, because if the theory is solved in the large N_c limit, one could study a perturbative expansion with coupling $1/N_c = 1/3$ and perhaps learn about the non-perturbative regime of QCD. In the course of these studies 't Hooft noticed that when $1/N_c$ is interpreted as a coupling strength, the resulting Feynman graph expansion is topologically identical to the worldsheet genus expansion of a generic interacting string theory. This was one of the early indications that Yang-Mills theory could be realized, in certain respects, as a theory of string.

In 1997 Maldacena fused 't Hooft's holographic principle and the $1/N_c$ expansion in a dramatic new proposal [20]. It was known that one can construct a four-dimensional maximally supersymmetric ($\mathcal{N} = 4$) $SU(N_c)$ gauge theory by stacking N_c coincident D3-branes and allowing open strings to stretch between pairs of branes [21]. The 't Hooft limit becomes accessible in this setting by taking the number of branes to be large. Since the D-branes are massive, however, a large number of them warp the ten-dimensional background geometry and a horizon is formed. The geometry in the near-horizon limit can be computed to be the product space of a five-dimensional anti-de-Sitter manifold and a five-dimensional sphere, or $AdS_5 \times S^5$. Furthermore, the branes are sources for closed string states, and the physics in the region just exterior to the branes is described by type IIB closed superstring theory in an $AdS_5 \times S^5$ background geometry. According to holography, the theory on the horizon should correspond to the physics inside the horizon. Maldacena was thereby led to conclude that type IIB superstring theory on $AdS_5 \times S^5$ is *equivalent* to $\mathcal{N} = 4$ supersymmetric Yang-Mills theory with $SU(N_c)$ gauge group in four spacetime dimensions! The conjectured equivalence of these two theories is a holographic duality. The relationship turns out to be dual in the more traditional sense, insofar as the coupling strengths that govern perturbative expansions in each theory are inversely proportional: perturbative physics in one theory corresponds to a non-perturbative regime

in the dual theory. The power afforded by a conjectured duality, however, is sometimes tempered by the inability to directly verify the proposal. Generically, a direct verification would require specific knowledge of non-perturbative physics on at least one side of the duality.

0.3 The Penrose limit

It should be noted that there is a substantial body of evidence that stands in support of Maldacena's conjecture. Most notably, the string and gauge theories are both invariant under the same superconformal symmetry group: $PSU(2, 2|4)$. Apart from the satisfaction of achieving a proof of the conjecture, an exploration of the underlying details would be useful in its own right; a more detailed understanding of how the AdS/CFT correspondence is realized on the microscopic level would be extremely valuable. The primary obstructions to such a program have been the difficulty of computing the dimensions of non-BPS operators in the strong-coupling limit of the gauge theory, and the unsolved problem of string quantization in the presence of a curved, Ramond-Ramond (RR) background geometry. In February of 2002, Berenstein, Maldacena and Nastase (BMN) found a specific set of limits where these problems can, to some extent, be circumvented [22]. In this section we will briefly review how this is achieved, paying particular attention to the string side of the duality (relevant details of the gauge theory will be covered in Chapter 1).

In convenient global coordinates, the $AdS_5 \times S^5$ metric can be written in the form

$$ds^2 = \widehat{R}^2 (-\cosh^2 \rho \, dt^2 + d\rho^2 + \sinh^2 \rho \, d\Omega_3^2 + \cos^2 \theta \, d\phi^2 + d\theta^2 + \sin^2 \theta \, d\widetilde{\Omega}_3^2) , \quad (0.3.1)$$

where \widehat{R} denotes the radius of both the sphere and the AdS space. (The hat is introduced because we reserve the symbol R for R -charge in the gauge theory.) The coordinate ϕ is periodic with period 2π and, strictly speaking, the time coordinate

t exhibits the same periodicity. In order to accommodate string dynamics, it is necessary to pass to the covering space in which time is *not* taken to be periodic. This geometry is accompanied by an RR field with N_c units of flux on the S^5 . It is a consistent, maximally supersymmetric type IIB superstring background provided that

$$\widehat{R}^4 = g_s N_c (\alpha')^2 , \quad (0.3.2)$$

where g_s is the string coupling. Explicitly, the AdS/CFT correspondence asserts that this string theory is equivalent to $\mathcal{N} = 4$ super Yang–Mills theory in four dimensions with an $SU(N_c)$ gauge group and coupling constant $g_{YM}^2 = g_s$. To simplify both sides of the correspondence, we study the duality in the simultaneous limits $g_s \rightarrow 0$ (the classical limit of the string theory) and $N_c \rightarrow \infty$ (the planar diagram limit of the gauge theory) with the 't Hooft coupling $g_{YM}^2 N_c$ held fixed. The holographically dual gauge theory is defined on the conformal boundary of $AdS_5 \times S^5$, which, in this case, is $R \times S^3$. Specifically, duality demands that operator dimensions in the conformally invariant gauge theory be equal to the energies of corresponding states of the ‘first-quantized’ string propagating in the $AdS_5 \times S^5$ background [23].

The quantization problem is simplified by boosting the string to lightlike momentum along some direction or, equivalently, by quantizing the string in the background obtained by taking a Penrose limit of the original geometry using the lightlike geodesic corresponding to the boosted trajectory. The simplest choice is to boost along an equator of the S^5 or, equivalently, to take a Penrose limit with respect to the lightlike geodesic $\phi = t$, $\rho = \theta = 0$. To perform lightcone quantization about this geodesic, it is helpful to make the reparameterizations

$$\cosh \rho = \frac{1 + z^2/4}{1 - z^2/4} , \quad \cos \theta = \frac{1 - y^2/4}{1 + y^2/4} , \quad (0.3.3)$$

and work with the metric

$$ds^2 = \widehat{R}^2 \left[- \left(\frac{1 + \frac{1}{4}z^2}{1 - \frac{1}{4}z^2} \right)^2 dt^2 + \left(\frac{1 - \frac{1}{4}y^2}{1 + \frac{1}{4}y^2} \right)^2 d\phi^2 + \frac{dz_k dz_k}{(1 - \frac{1}{4}z^2)^2} + \frac{dy_{k'} dy_{k'}}{(1 + \frac{1}{4}y^2)^2} \right], \quad (0.3.4)$$

where $y^2 = \sum_{k'} y^{k'} y^{k'}$ with $k' = 5, \dots, 8$ and $z^2 = \sum_k z^k z^k$ with $k = 1, \dots, 4$ define eight ‘Cartesian’ coordinates transverse to the geodesic. This metric is invariant under the full $SO(4, 2) \times SO(6)$ symmetry, but only translation invariance in t and ϕ and the $SO(4) \times SO(4)$ symmetry of the transverse coordinates remain manifest in this form. The translation symmetries mean that string states have a conserved energy ω , conjugate to t , and a conserved (integer) angular momentum J , conjugate to ϕ . Boosting along the equatorial geodesic is equivalent to studying states with large J and the lightcone Hamiltonian will give the (finite) allowed values for $\omega - J$ in that limit. On the gauge theory side, the S^5 geometry is replaced by an $SO(6)$ R -symmetry group, and J corresponds to the eigenvalue R of an $SO(2)$ R -symmetry generator. The AdS/CFT correspondence implies that string energies in the large- J limit should match operator dimensions in the limit of large R -charge.

On dimensional grounds, taking the $J \rightarrow \infty$ limit on string states is equivalent to taking the $\widehat{R} \rightarrow \infty$ limit of the geometry (in properly chosen coordinates). The coordinate redefinitions

$$t \rightarrow x^+, \quad \phi \rightarrow x^+ + \frac{x^-}{\widehat{R}^2}, \quad z_k \rightarrow \frac{z_k}{\widehat{R}}, \quad y_{k'} \rightarrow \frac{y_{k'}}{\widehat{R}} \quad (0.3.5)$$

make it possible to take a smooth $\widehat{R} \rightarrow \infty$ limit. (The lightcone coordinates x^\pm are a bit unusual, but have been chosen for future convenience in quantizing the worldsheet Hamiltonian.) Expressing the metric (0.3.4) in these new coordinates, we obtain the following expansion in powers of $1/\widehat{R}^2$:

$$ds^2 \approx 2 dx^+ dx^- + dz^2 + dy^2 - (z^2 + y^2) (dx^+)^2 + O(1/\widehat{R}^2). \quad (0.3.6)$$

The leading contribution (which we will call ds_{pp}^2) is the Penrose limit, or pp-wave geometry: it describes the geometry seen by the infinitely boosted string. The x^+ coordinate is dimensionless, x^- has dimensions of length squared, and the transverse coordinates now have dimensions of length.

In lightcone gauge quantization of the string dynamics, one identifies worldsheet time τ with the x^+ coordinate, so that the worldsheet Hamiltonian corresponds to the conjugate space-time momentum $p_+ = \omega - J$. Additionally, one sets the worldsheet momentum density $p_- = 1$ so that the other conserved quantity carried by the string, $p_- = J/\hat{R}^2$, is encoded in the length of the σ interval (though we will later keep p_- explicit for reasons covered in Chapter 3). Once x^\pm are eliminated, the quadratic dependence of ds_{pp}^2 on the remaining eight transverse bosonic coordinates leads to a quadratic (and hence soluble) bosonic lightcone Hamiltonian p_+ . Things are less simple when $1/\hat{R}^2$ corrections to the metric are taken into account: they add quartic interactions to the lightcone Hamiltonian and lead to nontrivial shifts in the spectrum of the string. This phenomenon, generalized to the superstring, will be the primary subject of this dissertation.

While it is clear how the Penrose limit can bring the bosonic dynamics of the string under perturbative control, the RR field strength survives this limit and causes problems for quantizing the superstring. The Green-Schwarz (GS) action is the only practical approach to quantizing the superstring in RR backgrounds, and we must construct this action for the IIB superstring in the $AdS_5 \times S^5$ background [24], pass to lightcone gauge and then take the Penrose limit. The latter step reduces the otherwise extremely complicated action to a worldsheet theory of free, equally massive transverse bosons and fermions [25]. As an introduction to the issues we will be concerned with, we give a concise summary of the construction and properties of the lightcone Hamiltonian $H_{\text{pp}}^{\text{GS}}$ that describe the superstring in this limit. This will be a helpful preliminary to our principal goal of evaluating the corrections to the Penrose limit of the GS action.

Gauge fixing eliminates the oscillating contributions to both lightcone coordinates

x^\pm , leaving eight transverse coordinates x^I as bosonic dynamical variables. Type IIB supergravity has two ten-dimensional supersymmetries that are described by two 16-component Majorana–Weyl spinors of the same ten-dimensional chirality. The GS superstring action contains just such a set of spinors (so that the desired spacetime supersymmetry comes out ‘naturally’). In the course of lightcone gauge fixing, half of these fermi fields are set to zero, leaving behind a complex eight-component worldsheet fermion ψ . This field is further subject to the condition that it transform in an $\mathbf{8}_s$ representation under $SO(8)$ rotations of the transverse coordinates (while the bosons of course transform as an $\mathbf{8}_v$). In a 16-component notation the restriction of the worldsheet fermions to the $\mathbf{8}_s$ representation is implemented by the condition $\gamma^9\psi = +\psi$ where $\gamma^9 = \gamma^1 \cdots \gamma^8$ and the γ^A are eight real, symmetric gamma matrices satisfying a Clifford algebra $\{\gamma^A, \gamma^B\} = 2\delta^{AB}$. Another quantity, which proves to be important in what follows, is $\Pi \equiv \gamma^1\gamma^2\gamma^3\gamma^4$. One could also define $\tilde{\Pi} = \gamma^5\gamma^6\gamma^7\gamma^8$, but $\Pi\psi = \tilde{\Pi}\psi$ for an $\mathbf{8}_s$ spinor.

In the Penrose limit, the lightcone GS superstring action takes the form

$$S_{\text{pp}} = \frac{1}{2\pi\alpha'} \int d\tau \int d\sigma (\mathcal{L}_B + \mathcal{L}_F) , \quad (0.3.7)$$

where

$$\mathcal{L}_B = \frac{1}{2} [(\dot{x}^A)^2 - (x'^A)^2 - (x^A)^2] , \quad (0.3.8)$$

$$\mathcal{L}_F = i\psi^\dagger \dot{\psi} + \psi^\dagger \Pi \psi + \frac{i}{2} (\psi \psi' + \psi^\dagger \psi'^\dagger) . \quad (0.3.9)$$

The fermion mass term $\psi^\dagger \Pi \psi$ arises from the coupling to the background RR five-form field strength, and matches the bosonic mass term (as required by supersymmetry). It is important that the quantization procedure preserve supersymmetry. However, as is typical in lightcone quantization, some of the conserved generators are linearly realized on the x^A and ψ^α , and others have a more complicated non-linear realization.

The equation of motion of the transverse string coordinates is

$$\ddot{x}^A - x''^A + x^A = 0 . \quad (0.3.10)$$

The requirement that x^A be periodic in the worldsheet coordinate σ (with period $2\pi\alpha'p_-$) leads to the mode expansion

$$x^A(\sigma, \tau) = \sum_{n=-\infty}^{\infty} x_n^A(\tau) e^{-ik_n\sigma} , \quad k_n = \frac{n}{\alpha'p_-} = \frac{n\widehat{R}^2}{\alpha'J} . \quad (0.3.11)$$

The canonical momentum p^A also has a mode expansion, related to that of x^A by the free-field equation $p^A = \dot{x}^A$. The coefficient functions are most conveniently expressed in terms of harmonic oscillator raising and lowering operators:

$$x_n^A(\tau) = \frac{i}{\sqrt{2\omega_n p_-}} (a_n^A e^{-i\omega_n\tau} - a_{-n}^{A\dagger} e^{i\omega_n\tau}) , \quad (0.3.12)$$

$$p_n^A(\tau) = \sqrt{\frac{\omega_n}{2p_-}} (a_n^A e^{-i\omega_n\tau} + a_{-n}^{A\dagger} e^{i\omega_n\tau}) . \quad (0.3.13)$$

The harmonic oscillator frequencies are determined by the equation of motion (0.3.10) to be

$$\omega_n = \sqrt{1 + k_n^2} = \sqrt{1 + (n\widehat{R}^2/\alpha'J)^2} = \sqrt{1 + (g_{YM}^2 N_c n^2/J^2)} , \quad (0.3.14)$$

where the mode index n runs from $-\infty$ to $+\infty$. (Because of the mass term, there is no separation into right-movers and left-movers.) The canonical commutation relations are satisfied by imposing the usual creation and annihilation operator algebra:

$$[a_m^A, a_n^{B\dagger}] = \delta_{mn} \delta^{AB} \Rightarrow [x^A(\sigma), p^B(\sigma')] = i2\pi\alpha' \delta(\sigma - \sigma') \delta^{AB} . \quad (0.3.15)$$

The fermion equation of motion is

$$i(\dot{\psi} + \psi'^{\dagger}) + \Pi\psi = 0 . \quad (0.3.16)$$

The expansion of ψ in terms of creation and annihilation operators is achieved by expanding the field in worldsheet momentum eigenstates

$$\psi(\sigma, \tau) = \sum_{n=-\infty}^{\infty} \psi_n(\tau) e^{-ik_n \sigma} , \quad (0.3.17)$$

which are further expanded in terms of convenient positive and negative frequency solutions of the fermion equation of motion:

$$\psi_n(\tau) = \frac{1}{\sqrt{4p_- \omega_n}} (e^{-i\omega_n \tau} (\Pi + \omega_n - k_n) b_n + e^{i\omega_n \tau} (1 - (\omega_n - k_n) \Pi) b_n^\dagger) . \quad (0.3.18)$$

The frequencies and momenta in this expansion are equivalent to those of the bosonic coordinates. In order to reproduce the anticommutation relations

$$\{\psi(\tau, \sigma), \psi^\dagger(\tau, \sigma')\} = 2\pi \alpha' \delta(\sigma - \sigma') , \quad (0.3.19)$$

we impose the standard oscillator algebra

$$\{b_m^\alpha, b_n^{\beta\dagger}\} = \frac{1}{2} (1 + \gamma_9)^{\alpha\beta} \delta_{m,n} . \quad (0.3.20)$$

The spinor fields ψ carry 16 components, but the $\mathbf{8}_s$ projection reduces this to eight anticommuting oscillators, exactly matching the eight transverse oscillators in the bosonic sector. The final expression for the lightcone Hamiltonian is

$$H_{\text{pp}}^{\text{GS}} = \sum_{n=-\infty}^{+\infty} \omega_n \left(\sum_A (a_n^A)^\dagger a_n^A + \sum_\alpha (b_n^\alpha)^\dagger b_n^\alpha \right) . \quad (0.3.21)$$

The harmonic oscillator zero-point energies nicely cancel between bosons and fermions for each mode n . The frequencies ω_n depend on the single parameter

$$\lambda' = g_{YM}^2 N_c / J^2 , \quad \omega_n = \sqrt{1 + \lambda' n^2} , \quad (0.3.22)$$

so that one can take J and $g_{YM}^2 N_c$ to be simultaneously large while keeping λ' fixed.

If λ' is kept fixed and small, ω_n may be expanded in powers of λ' , suggesting that contact with perturbative Yang–Mills gauge theory is possible.

The spectrum is generated by $8 + 8$ transverse oscillators acting on ground states labeled by an $SO(2)$ angular momentum taking integer values $-\infty < J < \infty$ (note that the oscillators themselves carry zero $SO(2)$ charge). Any combination of oscillators may be applied to a ground state, subject to the constraint that the sum of the oscillator mode numbers must vanish (this is the level-matching constraint, the only constraint not eliminated by lightcone gauge-fixing). The energies of these states are the sum of the individual oscillator energies (0.3.14), and the spectrum is very degenerate.¹ For example, the 256 states of the form $A_n^\dagger B_{-n}^\dagger |J\rangle$ for a given mode number n (where A^\dagger and B^\dagger each can be any of the $8+8$ bosonic and fermionic oscillators) all have the energy

$$p_+ = \omega - J = 2\sqrt{1 + (g_{YM}^2 N_c n^2 / J^2)} \sim 2 + (g_{YM}^2 N_c n^2 / J^2) + \dots \quad (0.3.23)$$

In the weak coupling limit ($\lambda' \rightarrow 0$) the degeneracy is even larger because the dependence on the oscillator mode number n goes away! This actually makes sense from the dual gauge theory point of view where $p_+ \rightarrow D - R$ (D is the dimension and R is the R -charge carried by gauge-invariant operators of large R): at zero coupling, operators have integer dimensions and the number of operators with $D - R = 2$, for example, grows with R , providing a basis on which string multiplicities are reproduced. Even more remarkably, BMN were able to show [22] that subleading terms in a λ' expansion of the string energies match the first perturbative corrections to the gauge theory operator dimensions in the large R -charge limit. We will further review the details of this agreement in Chapters 1 and 3.

More generally, we expect exact string energies in the $AdS_5 \times S^5$ background to have a joint expansion in the parameters λ' , defined above, and $1/J$. We also expect

¹Note that the $n = 0$ oscillators raise and lower the string energy by a protected amount $\delta p_+ = 1$, independent of the variable parameters. These oscillators play a special role, enlarging the degeneracy of the string states in a crucial way, and we will call them ‘zero-modes’ for short.

the degeneracies found in the $J \rightarrow \infty$ limit (for fixed λ') to be lifted by interaction terms that arise in the worldsheet Hamiltonian describing string physics at large but finite J . Large degeneracies must nevertheless remain in order for the spectrum to be consistent with the $PSU(2, 2|4)$ global supergroup that should characterize the exact string dynamics. The specific pattern of degeneracies should also match that of operator dimensions in the $\mathcal{N} = 4$ super Yang–Mills theory. Since the dimensions must be organized by the $PSU(2, 2|4)$ superconformal symmetry of the gauge theory, consistency is at least possible, if not guaranteed.

0.4 The $1/J$ expansion and post-BMN physics

As noted above, the matching achieved by BMN should not be confined to the Penrose (or large-radius) limit of the bulk theory, or to the large R -charge limit of the CFT. When the Penrose limit is lifted, finite-radius curvature corrections to the pp-wave geometry can be viewed as interaction perturbations to the free string theory, which, in turn, correspond to first-order corrections, in inverse powers of the R -charge, to the spectrum of anomalous dimensions in the gauge theory. With the hope that the underlying structure of the duality can be understood more clearly in this perturbative context, this dissertation is dedicated to exploring the AdS/CFT correspondence when these effects are included. In this section we will briefly review the work appearing in the literature upon which this thesis is based. In addition, we will also point out some of the more important developments that have appeared as part of the large body of research that has appeared following the original BMN paper.

In references [26] and [27], it was demonstrated that the first-order curvature corrections to the pp-wave superstring theory precisely reproduce finite R -charge corrections to the anomalous dimensions of so-called BMN operators, and exhibit the full $\mathcal{N} = 4$ extended supermultiplet structure of the dual gauge theory. The leading-order correction to the string theory gives rise to a complicated interacting theory of bosons and fermions in a curved RR background. While the steps taken

to quantize the resulting theory were fairly elaborate, it was demonstrated that they comprise a practical and correct method for defining the GS superstring action in that background. A detailed prescription for matching string states to gauge theory operators was given specifically in [27], along with a description of the procedure used to quantize the fully supersymmetric string theory and manage the set of second-class fermionic constraints that arise in lightcone gauge.

While the conjectured equivalence of the two theories emerged in this perturbative context in a remarkable manner, these studies also took advantage of the underlying duality structure of the correspondence. In particular, finite R -charge corrections to operator dimensions in the gauge theory emerge at all orders in $1/R$ (where R denotes the R -charge), but are defined perturbatively in the 't Hooft coupling $\lambda = g_{YM}^2 N$. Conversely, finite-radius corrections to string state energies appear perturbatively in inverse powers of the radius, or, equivalently, in inverse powers of the angular momentum J about the S^5 (which is identified with the gauge theory R -charge). According to duality, however, the string theory should provide a strong-coupling description of the gauge theory. This is realized by the fact that string energy corrections can be computed to all orders in the so-called modified 't-Hooft coupling $\lambda' = g_{YM}^2 N/J^2$. By studying the dilatation generator of $\mathcal{N} = 4$ SYM theory, several groups have been able to compute gauge theory operator dimensions to higher loop-order in λ (see, e.g., [28–36]), and, by expanding the corresponding string energy formulas in small λ' , the one- and two-loop energy corrections can be shown to precisely match the gauge theory results in a highly nontrivial way. The three-loop terms disagree, however, and this mismatch comprises a longstanding puzzle in these studies. Some investigations indicate that an order-of-limits issue may be responsible for this disagreement, whereby the small- λ expansion in the gauge theory fails to capture certain mixing interactions (known as wrapping terms) that are mediated by the dilatation generator [37].

To explore the correspondence further, and perhaps to shed light on the established three-loop disagreement, a complete treatment of the 4,096-dimensional space

of three-excitation string states was given in reference [38], including a comparison with corresponding SYM operators carrying three R -charge impurities. (The investigations in references [26] and [27] were restricted to the 256-dimensional space of two-excitation string states, also known as two-impurity states.) Although the interacting theory in this larger space is much more complicated, it was found that the full $\mathcal{N} = 4$ SYM extended supermultiplet structure is again realized by the string theory, and precise agreement with the anomalous dimension spectrum in the gauge theory was obtained to two-loop order in λ' . Once again, however, the three-loop formulas disagree.

Concurrent with these studies, a new formalism emerged for computing operator dimensions in the gauge theory. This began when Minahan and Zarembo were able to identify the one-loop mixing matrix of SYM operator dimensions with the Hamiltonian of an integrable $SO(6)$ spin chain with vector lattice sites [39]. One practical consequence of this discovery is that the quantum spin chain Hamiltonian describing the SYM dilatation generator can be completely diagonalized by a set of algebraic relations known as the Bethe ansatz. Work in the $SO(6)$ sector was extended by Beisert and Staudacher, who formulated a Bethe ansatz for the full $PSU(2, 2|4)$ superconformal symmetry of the theory (under which the complete dilatation generator is invariant) [32].

The emergence of integrable structures in the gauge theory has given rise to many novel tests of AdS/CFT (see, e.g., [40–58]). It has been suggested by Bena, Polchinski and Roiban, for instance, that the classical lightcone gauge worldsheet action of type IIB superstring theory in $AdS_5 \times S^5$ may itself be integrable [59]. If both theories are indeed integrable, they should admit infinite towers of hidden charges that, in turn, should be equated via the AdS/CFT correspondence, analogous to identifying the SYM dilatation generator with the string Hamiltonian. Numerous investigations have been successful in matching classically conserved hidden string charges with corresponding charges derived from the integrable structure of the gauge theory. Arutyunov and Staudacher, for example, were able to show that an infinite set of local

charges generated via Bäcklund transformations on certain classical extended string solutions can be matched to an infinite tower of charges generated by a corresponding sector of gauge theory operators [41]. It is important to note, however, that these identifications are between the structures of *classically* integrable string sigma models and integrable *quantum* spin chains. Along these lines of investigation, Arutyunov, Frolov and Staudacher developed an interpolation between the classical string sigma model and the quantum spin chain that yielded a Bethe ansatz purported to capture the dynamics of an $SU(2)$ sector of the string theory [44]. This ansatz, though conjectural, allowed the authors to extract multi-impurity string energy predictions in the near-pp-wave limit (at $O(1/J)$ in the curvature expansion). Corresponding predictions were extracted in reference [60] directly from the quantized string theory, and the resulting formulas matched the Bethe ansatz predictions to all loop-orders in λ' in a remarkable and highly intricate fashion.

Recently the question of quantum integrability in the string theory was addressed in reference [61]. Using a perturbed Lax representation of a particular solitonic solution to the string sigma model, one is able to argue that the string theory admits an infinite tower of hidden commuting charges that are conserved by the quantized theory to quartic order in field fluctuations. In addition, a prescription for matching the eigenvalue spectra of these charges to dual quantities in the gauge theory can also be formulated.

At this point there is a considerable amount of evidence that both the string and gauge theories are exactly integrable (see also [62, 63] for recent developments). The hope is of course that we will ultimately be led to an exact solution to large- N_c Yang-Mills theory. Before reaching this goal, it is reasonable to expect that type IIB string theory on $AdS_5 \times S^5$ and $\mathcal{N} = 4$ super Yang-Mills theory will be shown to admit identical Bethe ansatz equations, thereby proving this particular duality. This is likely the next major step in these investigations. There are several intermediate problems that need to be solved, however, including the known mismatch between the string and gauge theory at three-loop order in the 't Hooft coupling. The resolution

of these outstanding problems will inevitably lead to a deeper understanding of both the relationship between gauge and string theory, and the capacity of string theory itself to generate realistic models of particle physics.

0.5 Overview

In this dissertation we will work in the large- N_c limit, where we can ignore string splitting and joining interactions; the “stringy” effects we are concerned with arise strictly from interactions among the bosonic and fermionic field excitations on the worldsheet. In Chapter 1 we will provide a brief treatment of the relevant calculations that are needed on the gauge theory side of the correspondence, based on work originally presented in [26]. While the results computed there can be found elsewhere in the literature (see, e.g., [28]), we present our own derivation for pedagogical reasons and to arrange the computation in a way that clarifies the eventual comparison with string theory.

As noted above, the task of calculating operator dimensions in the planar limit of $\mathcal{N} = 4$ super Yang-Mills theory can be vastly simplified by mapping the dilation generator to the Hamiltonian of an integrable spin chain. These techniques are powerful at leading order in perturbation theory but become increasingly complicated beyond one loop in the ’t Hooft parameter $\lambda = g_{\text{YM}}^2 N_c$, where spin chains typically acquire long-range (non-nearest-neighbor) interactions. In certain sectors of the theory, moreover, higher-loop Bethe ansätze do not even exist. In Chapter 2 we develop a virial expansion of the spin chain Hamiltonian as an alternative to the Bethe ansatz methodology, a method that simplifies the computation of dimensions of multi-impurity operators at higher loops in λ . We use these methods to extract numerical gauge theory predictions near the BMN limit for comparison with corresponding results on the string theory side of the AdS/CFT correspondence. For completeness, we compare our virial results with predictions that can be derived from current Bethe ansatz technology.

In Chapter 3 we compute the complete set of first curvature corrections to the lightcone gauge string theory Hamiltonian that arise in the expansion of $AdS_5 \times S^5$ about the pp-wave limit. We develop a systematic quantization of the interacting worldsheet string theory and use it to obtain the interacting spectrum of the so-called ‘two-impurity’ states of the string. The quantization is technically rather intricate and we provide a detailed account of the methods we use to extract explicit results. We give a systematic treatment of the fermionic states and are able to show that the spectrum possesses the proper extended supermultiplet structure (a nontrivial fact since half the supersymmetry is nonlinearly realized). We test holography by comparing the string energy spectrum with the scaling dimensions of corresponding gauge theory operators. We show that agreement is obtained in low orders of perturbation theory, but breaks down at third order.

Notwithstanding this third-order mismatch, we proceed with this line of investigation in Chapter 4 by subjecting the string and gauge theories to significantly more rigorous tests. Specifically, we extend the results of Chapter 3 at $O(1/J)$ in the curvature expansion to include string states and SYM operators with three worldsheet or R -charge impurities. In accordance with the two-impurity problem, we find a perfect and intricate agreement between both sides of the correspondence to two-loop order in λ and, once again, the string and gauge theory predictions fail to agree at third order.

In Chapter 5 we generalize this analysis on the string side by directly computing string energy eigenvalues in certain protected sectors of the theory for an arbitrary number of worldsheet excitations with arbitrary mode-number assignments. While our results match all existing gauge theory predictions to two-loop order in λ' , we again observe a mismatch at three loops between string and gauge theory. We find remarkable agreement to *all* loops in λ' , however, with the near pp-wave limit of a Bethe ansatz for the quantized string Hamiltonian given in an $\mathfrak{su}(2)$ sector. Based on earlier two- and three-impurity results, we also infer the full multiplet decomposition of the N -impurity superstring theory with distinct mode excitations to two loops in

λ' .

In Chapter 6 we build on recent explorations of the AdS/CFT correspondence that have unveiled integrable structures underlying both the gauge and string theory sides of the correspondence. By studying a semiclassical expansion about a class of point-like solitonic solutions to the classical string equations of motion on $AdS_5 \times S^5$, we take a step toward demonstrating that integrability in the string theory survives quantum corrections beyond tree level. Quantum fluctuations are chosen to align with background curvature corrections to the pp-wave limit of $AdS_5 \times S^5$, and we present evidence for an infinite tower of local bosonic charges that are conserved by the quantum theory to quartic order in the expansion. We explicitly compute several higher charges based on a Lax representation of the worldsheet sigma model and provide a prescription for matching the eigenvalue spectra of these charges with corresponding quantities descending from the integrable structure of the gauge theory. The final chapter is dedicated to a discussion of the current status of these studies and an overview of future directions of investigation.

Chapter 1

$\mathcal{N} = 4$ super Yang-Mills theory

As discussed in the introduction, the AdS/CFT correspondence states that the energy spectrum of string excitations in an anti-de-Sitter background should be equivalent (albeit related by a strong/weak duality) to the spectrum of operator anomalous dimensions of the field theory living on the conformal boundary of that background. Any attempt to test the validity of this statement directly must therefore involve a computation of operator dimensions in the gauge theory, particularly for those operators that are non-BPS. As discussed above, this is a nontrivial task for generic gauge theory operators, but the advent of the BMN mechanism has led to dramatic simplifications and insights. Following the appearance of the original BMN paper [22], the field witnessed remarkable progress in understanding the dilatation generator of $\mathcal{N} = 4$ SYM theory (see, e.g., [28–36, 40–58]). The review presented in this chapter will focus on some of the major contributions to this understanding. Since this work is dedicated primarily to understanding the string theory side of the AdS/CFT correspondence, special preference will be given to information that contributes directly to our ability to interpret the dual spectrum of string excitations. For a more comprehensive and detailed review of the gauge theory aspects of these studies, the reader is referred to [35].

To arrange the calculation in a way that is more useful for our subsequent comparison with string theory, and to emphasize a few specific points, it is useful to rederive several important results. We will focus in Section 1.1 on the dimensions

and multiplicities of a specific set of near-BPS (two-impurity) operators in the planar limit. Most of the information to be covered in this section originally appeared in [28], though we will orient our review around a rederivation of these results first presented in [26]. Section 1.2 generalizes these results to the complete set of two-impurity, single-trace operators. This will set the stage for a detailed analysis of the corresponding string energy spectrum.

1.1 Dimensions and multiplicities

As explained above, the planar large- N_c limit of the gauge theory corresponds to the noninteracting sector ($g_s \rightarrow 0$) of the dual string theory.¹ In this limit the gauge theory operators are single-trace field monomials classified by dimension D and the scalar $U(1)_R$ component (denoted by R) of the $SU(4)$ R -symmetry group. We will focus in this section on the simple case of operators containing only two R -charge impurities. The classical dimension will be denoted by K , and the BMN limit is reached by taking $K, R \rightarrow \infty$ such that $\Delta_0 \equiv K - R$ is a fixed, finite integer. The anomalous dimensions (or $D - K$) are assumed to be finite in this limit, and the quantity $\Delta \equiv D - R$ is defined for comparison with the string lightcone Hamiltonian $P_+ = \omega - J$ (see Section 0.3 of the introduction).

It is useful to classify operators in the gauge theory according to their representation under the exact global $SU(4)$ R -symmetry group. This is possible because the dimension operator commutes with the R -symmetry. We therefore find it convenient to label the component fields with Young boxes, which clarifies the decomposition of composite operators into irreducible tensor representations of $SU(4)$. More specifically, the tensor irreps of $SU(4)$ are represented by Young diagrams composed of at most three rows of boxes denoted by a set of three numbers (n_1, n_2, n_3) indicating the differences in length of successive rows. The fields available are a gauge field, a set of gluinos transforming as $\mathbf{4}$ and $\bar{\mathbf{4}}$ under the R -symmetry group, and a set of

¹The Yang-Mills genus-counting parameter is $g_2 = J^2/N_c$ [64, 65].

scalars transforming as a **6**. In terms of Young diagrams, the gluinos transform as two-component Weyl spinors in the $(1,0,0)$ fundamental (**4**) and its adjoint $(0,0,1)$ in the antifundamental ($\bar{\mathbf{4}}$):

$$\chi_a^\square \text{ (**4**)}, \quad \bar{\chi}_{\dot{a}}^{\square} \text{ ($\bar{\mathbf{4}}$)}.$$

The a and \dot{a} indices denote transformation in the $(\mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2})$ representations of $SL(2, C)$ (the covering group of the spacetime Lorentz group), respectively. Likewise, the scalars appear as

$$\phi^{\square} \text{ (**6**)}.$$

In the planar large- N_c limit the operators of interest are those containing only a single gauge trace. To work through an explicit example, we will restrict attention for the moment to operators comprising spacetime scalars. It is convenient to further classify these operators under the decomposition

$$SU(4) \supset SU(2) \times SU(2) \times U(1)_R, \quad (1.1.1)$$

since we are eventually interested in taking the scalar $U(1)_R$ component to be large (which corresponds to the large angular momentum limit of the string theory). The $U(1)_R$ charge of the component fields above can be determined by labeling the Young diagrams attached to each field with $SU(4)$ indices, assigning $R = \frac{1}{2}$ to the indices 1, 2 and $R = -\frac{1}{2}$ to the indices 3, 4:

$$\begin{aligned} R = 1 : \phi^{\begin{smallmatrix} \square \\ 2 \end{smallmatrix}}(Z), \quad R = 0 : \phi^{\begin{smallmatrix} \square \\ 3 \end{smallmatrix}}, \phi^{\begin{smallmatrix} \square \\ 4 \end{smallmatrix}}, \phi^{\begin{smallmatrix} \square \\ 3 \end{smallmatrix}}, \phi^{\begin{smallmatrix} \square \\ 4 \end{smallmatrix}}(\phi^A), \quad R = -1 : \phi^{\begin{smallmatrix} \square \\ 4 \end{smallmatrix}}(\bar{Z}), \\ R = 1/2 : \chi^{\begin{smallmatrix} \square \\ 1 \end{smallmatrix}}, \chi^{\begin{smallmatrix} \square \\ 2 \end{smallmatrix}}, \bar{\chi}^{\begin{smallmatrix} \square \\ 3 \end{smallmatrix}}, \bar{\chi}^{\begin{smallmatrix} \square \\ 4 \end{smallmatrix}}, \quad R = -1/2 : \chi^{\begin{smallmatrix} \square \\ 3 \end{smallmatrix}}, \chi^{\begin{smallmatrix} \square \\ 4 \end{smallmatrix}}, \bar{\chi}^{\begin{smallmatrix} \square \\ 3 \end{smallmatrix}}, \bar{\chi}^{\begin{smallmatrix} \square \\ 4 \end{smallmatrix}}. \end{aligned} \quad (1.1.2)$$

To remain consistent with the literature we have labeled the scalars using either Z

or \bar{Z} for fields with $R = 1$ or $R = -1$, respectively, or ϕ^A (with $A \in 1, \dots, 4$) for fields with zero R -charge. The types of operators of interest to us are those with large naive dimension K and large R -charge, with the quantity $\Delta_0 \equiv K - R$ held fixed. The number Δ_0 is typically referred to as the impurity number of the operator; as explained above, N -impurity SYM operators map to string states created by N oscillators acting on the vacuum, subject to level matching. Operators in the gauge theory with zero impurity number are BPS, and their dimensions are protected. The first interesting set of non-BPS operators are those with $\Delta_0 = 2$. Restricting to spacetime scalars with $\Delta_0 \leq 2$, we have

$$\begin{aligned}
& \text{tr}((\phi^{\square})^K), & (R_{\max} = K) \\
& \text{tr}((\chi^{\square}\sigma_2\chi^{\square})(\phi^{\square})^{K-3}), \quad \text{tr}((\chi^{\square}\phi^{\square}\sigma_2\chi^{\square})(\phi^{\square})^{K-4}), \quad \dots & (R_{\max} = K - 2) \\
& \text{tr}((\bar{\chi}^{\square}\sigma_2\bar{\chi}^{\square})(\phi^{\square})^{K-3}), \quad \text{tr}((\bar{\chi}^{\square}\phi^{\square}\sigma_2\bar{\chi}^{\square})(\phi^{\square})^{K-4}), \quad \dots & (R_{\max} = K - 2) \\
& \text{tr}(\nabla_{\mu}\phi^{\square}\nabla^{\mu}\phi^{\square}(\phi^{\square})^{K-4}), & (R_{\max} = K - 2),
\end{aligned} \tag{1.1.3}$$

where ∇ is the spacetime gauge-covariant derivative.

Starting with purely bosonic operators with no derivative insertions, we must decompose into irreps an $SU(4)$ tensor of rank $2K$. These irreps are encoded in Young diagrams with $2K$ total boxes, and the goal is to determine the multiplicity with which each diagram appears. (An alternative approach, taken in [28], is to use the bosonic $SO(6)$ sector of the R -symmetry group.) For the purposes of this example, we restrict to irreducible tensors in the expansion with $\Delta_0 = 0, 2$. For K odd we have

$$\begin{aligned}
\text{tr}(\phi^{\square}{}^K) \rightarrow & 1 \times \underbrace{\begin{array}{|c|c|c|c|c|} \hline \square & \square & \square & \square & \square \\ \hline \end{array}}_K \oplus \left(\frac{K-1}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|} \hline \square & \square & \square & \square \\ \hline \end{array}}_{K-1} \oplus \left(\frac{K-1}{2}\right) \times \underbrace{\begin{array}{|c|c|c|} \hline \square & \square & \square \\ \hline \end{array}}_{K-2} \\
& \oplus \left(\frac{K-1}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|} \hline \square & \square & \square & \square \\ \hline \end{array}}_{K-1} \oplus \left(\frac{K-3}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|c|} \hline \square & \square & \square & \square & \square \\ \hline \end{array}}_K \oplus \dots, \tag{1.1.4}
\end{aligned}$$

while for K even we have

$$\begin{aligned} \text{tr}(\phi^{\square K}) \rightarrow & 1 \times \underbrace{\square}_{K} \oplus \left(\frac{K-2}{2}\right) \times \underbrace{\square}_{K-1} \oplus \left(\frac{K}{2}\right) \times \underbrace{\square}_{K-2} \\ & \oplus \left(\frac{K-2}{2}\right) \times \underbrace{\square}_{K-1} \oplus \left(\frac{K-2}{2}\right) \times \underbrace{\square}_{K} \oplus \dots \end{aligned} \quad (1.1.5)$$

The irreps with larger minimal values of $\Delta_0 = K - R$ have multiplicities that grow as higher powers of K . This is very significant for the eventual string theory interpretation of the anomalous dimensions, but we will not expand on this point here.

The bifermion operators (that are spacetime scalars) with $\Delta_0 = 2$ contain products of two gluinos and $K - 3$ scalars:

$$\text{tr}(\chi^{\square} \sigma_2 \chi^{\square} (\phi^{\square})^{K-3}) \rightarrow 1 \times \underbrace{\square}_{K-2} \oplus 1 \times \underbrace{\square}_{K-1} \oplus \dots, \quad (1.1.6)$$

$$\text{tr}(\bar{\chi}^{\square} \sigma_2 \bar{\chi}^{\square} (\phi^{\square})^{K-3}) \rightarrow 1 \times \underbrace{\square}_{K-1} \oplus 1 \times \underbrace{\square}_{K-2} \oplus \dots \quad (1.1.7)$$

Note that products of χ^{\square} and $\bar{\chi}^{\square}$ cannot be made to form spacetime scalars because they transform under inequivalent irreps of $SL(2, C)$.

Different operators are obtained by different orderings of the component fields, but such operators are not necessarily independent under cyclic permutations or permutations of the individual fields themselves, subject to the appropriate statistics. Using an obvious shorthand notation, the total multiplicities of bifermion irreps are as follows for K odd:

$$\text{tr}(\chi^{\square} \sigma_2 \chi^{\square} (\phi^{\square})^{K-3}) \rightarrow \left(\frac{K-3}{2}\right) \times \underbrace{\square}_{K-2} \oplus \left(\frac{K-1}{2}\right) \times \underbrace{\square}_{K-1} \oplus \dots, \quad (1.1.8)$$

$$\text{tr}(\bar{\chi}^{\square} \sigma_2 \bar{\chi}^{\square} (\phi^{\square})^{K-3}) \rightarrow \left(\frac{K-3}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|} \hline \square & \square & \square & \square \\ \hline \square & \square & \square & \square \\ \hline \end{array}}_{K-2} \oplus \left(\frac{K-1}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|c|} \hline \square & \square & \square & \square & \square \\ \hline \square & \square & \square & \square & \square \\ \hline \square & \square & \square & \square & \square \\ \hline \end{array}}_{K-1} \oplus \dots \quad (1.1.9)$$

The results for K even are, once again, slightly different:

$$\text{tr}(\chi^{\square} \sigma_2 \chi^{\square} (\phi^{\square})^{K-3}) \rightarrow \left(\frac{K-2}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|} \hline \square & \square & \square & \square \\ \hline \square & \square & \square & \square \\ \hline \end{array}}_{K-2} \oplus \left(\frac{K-2}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|} \hline \square & \square & \square & \square \\ \hline \square & \square & \square & \square \\ \hline \end{array}}_{K-1} \oplus \dots, \quad (1.1.10)$$

$$\text{tr}(\bar{\chi}^{\square} \sigma_2 \bar{\chi}^{\square} (\phi^{\square})^{K-3}) \rightarrow \left(\frac{K-2}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|} \hline \square & \square & \square & \square \\ \hline \square & \square & \square & \square \\ \hline \end{array}}_{K-2} \oplus \left(\frac{K-2}{2}\right) \times \underbrace{\begin{array}{|c|c|c|c|c|} \hline \square & \square & \square & \square & \square \\ \hline \square & \square & \square & \square & \square \\ \hline \square & \square & \square & \square & \square \\ \hline \end{array}}_{K-1} \oplus \dots \quad (1.1.11)$$

Since the dimension operator can only have matrix elements between operators belonging to the same $SU(4)$ irrep, this decomposition amounts to a block diagonalization of the problem. The result of this program can be summarized by first noting that the decomposition can be divided into a BPS and non-BPS sector. The BPS states ($\Delta_0 = 0$) appear in the $(0, K, 0)$ irrep and do not mix with the remaining non-BPS sectors, which yield irreps whose multiplicities scale roughly as $K/2$ for large K . Even at this stage it is clear that certain irreps only appear in the decomposition of certain types of operators. The $(2, K-4, 2)$ irrep, for example, will only appear within the sector of purely bosonic operators (the same statement does not hold for the $(0, K-3, 2)$ irrep). Restricting to the $(2, K-4, 2)$ irrep, we see that the dimension matrix cannot mix operators in the purely bosonic sector with bifermions, for example. We will eventually make these sorts of observations much more precise, as they will become invaluable in subsequent analyses. The general problem involves diagonalizing matrices that are approximately $K/2 \times K/2$ in size. The operators of interest will have large $K = R + 2$ and fixed $\Delta_0 = K - R = 2$. As noted above, we expect that the anomalous dimension spectrum should match the energy spectrum of string states created by two oscillators acting on a ground state with angular

momentum $J = R$.

As an example we will start with the basis of $K - 1$ purely bosonic operators with dimension K and $\Delta_0 = 2$. The anomalous dimensions are the eigenvalues of the mixing matrix d_1^{ab} , appearing in the perturbative expansion of the generic two-point function according to

$$\langle O_a(x) O_b(0) \rangle \sim (x)^{-2d_0} (\delta_{ab} + \ln(x^2) d_1^{ab}) , \quad (1.1.12)$$

where d_0 is the naive dimension. The δ_{ab} term implies that the operator basis is orthonormal in the free theory (in the large- N_c limit, this is enforced by multiplying the operator basis by a common overall normalization constant). The operator basis can be expressed as

$$\{O_{K,1}^{AB}, \dots, O_{K,K-1}^{AB}\} = \{\text{tr}(ABZ^{K-2}), \text{tr}(AZBZ^{K-3}), \dots, \text{tr}(AZ^{K-3}BZ), \text{tr}(AZ^{K-2}B)\} , \quad (1.1.13)$$

where Z stands for $\phi^{\frac{0}{2}}$ and has $R = 1$, while A, B stand for any of the four ϕ^A ($A = 1, \dots, 4$) with $R = 0$ (the so-called R -charge impurities). The overall constant needed to orthonormalize this basis is easy to compute, but is not needed for the present purposes. Since the R -charge impurities A and B are $SO(4)$ vectors, the operators in this basis are rank-two $SO(4)$ tensors. In the language of $SO(4)$ irreps, the symmetric-traceless tensor descends from the $SU(4)$ irrep labeled by the $(2, K - 4, 2)$ Young diagram. Likewise, the antisymmetric tensor belongs to the pair $(0, K - 3, 2) + (2, K - 3, 0)$, and the $SO(4)$ trace (when completed to a full $SO(6)$ trace) belongs to the $(0, K - 2, 0)$ irrep. In what follows, we refer to these three classes of operator as $\overline{T}_K^{(+)}$, $\overline{T}_K^{(-)}$ and $\overline{T}_K^{(0)}$, respectively. If we take $A \neq B$, the trace part drops out and the $\overline{T}_K^{(\pm)}$ operators are isolated by symmetrizing and antisymmetrizing on A, B .

At one-loop order in the 't Hooft coupling $g_{YM}^2 N_c$ the action of the dilatation operator on the basis in eqn. (1.1.13), correct to all orders in $1/K$, produces a sum

of interchanges of all nearest-neighbor fields in the trace. All diagrams that exchange fields at greater separation (at this loop order) are non-planar, and are suppressed by powers of $1/N_c$. As an example, we may restrict to the $A \neq B$ case. Omitting the overall factor coming from the details of the Feynman diagram, the leading action of the anomalous dimension on the $K - 1$ bosonic monomials of (1.1.13) has the following structure:

$$\begin{aligned}
(ABZ^{K-2}) &\rightarrow (BAZ^{K-2}) + 2(AZBZ^{K-3}) + (K-3)(ABZ^{K-2}) , \\
(AZBZ^{K-3}) &\rightarrow 2(ABZ^{K-2}) + 2(AZ^2BZ^{K-4}) + (K-4)(AZBZ^{K-3}) , \\
&\dots\dots\dots \\
(AZ^{K-2}B) &\rightarrow 2(AZ^{K-3}BZ) + (K-3)(BAZ^{K-2}) + (ABZ^{K-2}) . \quad (1.1.14)
\end{aligned}$$

Arranging this into matrix form, we have

$$[\text{Anom Dim}]_{(K-1) \times (K-1)} \sim \begin{pmatrix} K-3 & 2 & 0 & \dots & 1 \\ 2 & K-4 & 2 & \dots & 0 \\ & & \ddots & & \\ 0 & \dots & 2 & K-4 & 2 \\ 1 & \dots & 0 & 2 & K-3 \end{pmatrix} . \quad (1.1.15)$$

As a final step, we must observe that the anomalous dimension matrix in eqn. (1.1.15) contains contributions from the $SU(4)$ irrep $(0, K, 0)$, which corresponds to the chiral primary $\text{tr}(Z^K)$. The eigenstate associated with this operator is $\vec{X}_0 = (1, \dots, 1)$, with eigenvalue K (the naive dimension). Since this operator is BPS, however, its anomalous dimension must be zero: to normalize the (1.1.15) we therefore subtract K times the identity, leaving

$$[\text{Anom Dim}]_{(K-1) \times (K-1)} \sim \begin{pmatrix} -3 & +2 & 0 & \dots & 1 \\ +2 & -4 & +2 & \dots & 0 \\ & & \ddots & & \\ 0 & \dots & +2 & -4 & +2 \\ +1 & \dots & 0 & +2 & -3 \end{pmatrix} . \quad (1.1.16)$$

The zero eigenvector belonging to the $(0, K, 0)$ representation should then be dropped. The anomalous dimensions are thus the nonzero eigenvalues of (1.1.16). This looks very much like the lattice Laplacian for a particle hopping from site to site on a periodic lattice. The special structure of the first and last rows assigns an extra energy to the particle when it hops past the origin. This breaks strict lattice translation invariance but makes sense as a picture of the dynamics involving two-impurity states: the impurities propagate freely when they are on different sites and have a contact interaction when they collide. This picture has led people to map the problem of finding operator dimensions onto the technically much simpler one of finding the spectrum of an equivalent quantum-mechanical Hamiltonian [66]; this important topic will be reserved for later chapters.

To determine the $SU(4)$ irrep assignment of each of the eigenvalues of (1.1.16), note that the set of operator monomials is invariant under $A \leftrightarrow B$. For some vector $\vec{C} = (C_1, \dots, C_{K-1})$ representing a given linear combination of monomials, this transformation sends $C_i \rightarrow C_{K-i}$. The matrix (1.1.16) itself is invariant under $A \leftrightarrow B$, so its eigenvectors will either be even ($C_i = C_{K-i}$) or odd ($C_i = -C_{K-i}$) under the same exchange. The two classes of eigenvalues and normalized eigenvectors are:

$$\begin{aligned} \lambda_n^{(K+)} &= 8 \sin^2 \left(\frac{n\pi}{K-1} \right), \quad n = 1, 2, \dots, n_{max} = \begin{cases} (K-3)/2 & K \text{ odd} \\ (K-2)/2 & K \text{ even} \end{cases}, \\ C_{n,i}^{(K+)} &= \frac{2}{\sqrt{K-1}} \cos \left[\frac{2\pi n}{K-1} \left(i - \frac{1}{2} \right) \right], \quad i = 1, \dots, K-1, \end{aligned} \quad (1.1.17)$$

$$\begin{aligned} \lambda_n^{(K-)} &= 8 \sin^2 \left(\frac{n\pi}{K} \right), \quad n = 1, 2, \dots, n_{max} = \begin{cases} (K-1)/2 & K \text{ odd} \\ (K-2)/2 & K \text{ even} \end{cases}, \\ C_{n,i}^{(K-)} &= \frac{2}{\sqrt{K}} \sin \left[\frac{2\pi n}{K} (i) \right], \quad i = 1, \dots, K-1. \end{aligned} \quad (1.1.18)$$

The eigenoperators are constructed from the eigenvectors according to

$$\bar{T}_{K,n}^{(\pm)}(x) = \sum_{i=1}^{K-1} C_{n,i}^{(K\pm)} O_{K,i}^{AB}(x) . \quad (1.1.19)$$

By appending the appropriate overall normalization factor and adding the zeroth order value $\Delta_0 = 2$, we obtain $\Delta = D - R$. The results are divided according to operators belonging to the $(2, K-4, 2)$ irrep $(\bar{T}_K^{(+)})$, the $(0, K-3, 2) + (2, K-3, 0)$ irreps $(\bar{T}_K^{(-)})$ and $(0, K-2, 0)$ $(\bar{T}_K^{(0)})$. In $SO(4)$ language, these are the symmetric-traceless, antisymmetric and trace representations, as described above. We therefore have the following, exact in K :

$$\begin{aligned} \Delta(\bar{T}_K^{(+)}) &= 2 + \frac{g_{YM}^2 N_c}{\pi^2} \sin^2 \left(\frac{n\pi}{K-1} \right) , \quad n = 1, 2, \dots, n_{max} = \begin{cases} (K-3)/2 & K \text{ odd} \\ (K-2)/2 & K \text{ even} \end{cases} , \\ \Delta(\bar{T}_K^{(-)}) &= 2 + \frac{g_{YM}^2 N_c}{\pi^2} \sin^2 \left(\frac{n\pi}{K} \right) , \quad n = 1, 2, \dots, n_{max} = \begin{cases} (K-1)/2 & K \text{ odd} \\ (K-2)/2 & K \text{ even} \end{cases} , \\ \Delta(\bar{T}_K^{(0)}) &= 2 + \frac{g_{YM}^2 N_c}{\pi^2} \sin^2 \left(\frac{n\pi}{K+1} \right) , \quad n = 1, 2, \dots, n_{max} = \begin{cases} (K-1)/2 & K \text{ odd} \\ (K/2) & K \text{ even} \end{cases} . \end{aligned} \quad (1.1.20)$$

The multiplicities match the earlier predictions given by the expansion in Young diagrams in eqns. (1.1.4) and (1.1.5).

We will eventually be interested in exploring the overlap of such results with that which can be predicted by the dual string theory. As described above, the central assumption introduced by Berenstein, Maldacena and Nastase is that the R -charge and the rank of the gauge group N_c can be taken to infinity such that the quantity N_c/R^2 remains fixed. The perturbation expansion in the gauge theory is then controlled by $g_{YM}^2 N_c$ (which is kept small in the $g_{YM}^2 \rightarrow 0$ limit, which is the

classical $g_s \rightarrow 0$ limit of the string theory), while worldsheet interactions in the string theory are controlled by $1/\widehat{R}$. If we express the dimension formulas (1.1.20) in terms of R -charge R , rather than naive dimension K (using $K = R + 2$) and take the limit in this way, we find

$$\begin{aligned}\Delta(\overline{T}_{R+2}^{(+)}) &\rightarrow 2 + \frac{g_{YM}^2 N_c}{R^2} n^2 \left(1 - \frac{2}{R} + O(R^{-2}) \right) , \\ \Delta(\overline{T}_{R+2}^{(-)}) &\rightarrow 2 + \frac{g_{YM}^2 N_c}{R^2} n^2 \left(1 - \frac{4}{R} + O(R^{-2}) \right) , \\ \Delta(\overline{T}_{R+2}^{(0)}) &\rightarrow 2 + \frac{g_{YM}^2 N_c}{R^2} n^2 \left(1 - \frac{6}{R} + O(R^{-2}) \right) .\end{aligned}\tag{1.1.21}$$

The key fact is that the degeneracy of the full BMN limit (at leading order in $1/R$) is lifted at subleading order in $1/R$. By including these subleading terms we generate an interesting spectrum that will prove to be a powerful tool for comparison with string theory and testing the claims of the AdS/CFT correspondence.

1.2 The complete supermultiplet

We have thus far reviewed the anomalous dimension computation for a specific set of operators. For a complete comparison with the string theory, we need to carry out some version of the above arguments for all the relevant operators with $\Delta_0 = 2$. While this is certainly possible, we can instead rely on supersymmetry to determine the full spectrum of anomalous dimensions for all single-trace, two-impurity operators. The extended superconformal symmetry of the gauge theory means that operator dimensions will be organized into multiplets based on a lowest-dimension primary \mathcal{O}_D of dimension D . Other conformal primaries within the multiplet can be generated by acting on super-primaries with any of eight supercharges that increment the anomalous dimension shifts by a fixed amount but leave the impurity number unchanged. We need only concern ourselves here with the case in which \mathcal{O}_D is a spacetime scalar (of dimension D and R -charge R). There are sixteen supercharges and we can choose

eight of them to be raising operators; there are $2^8 = 256$ operators we can reach by ‘raising’ the lowest one. Since the raising operators increase the dimension and R -charge by $1/2$ each time they act, the operators at level L , obtained by acting with L supercharges, all have the same dimension and R -charge. The corresponding decomposition of the 256-dimensional multiplet is shown in table 1.1.

Level	0	1	2	3	4	5	6	7	8
Multiplicity	1	8	28	56	70	56	28	8	1
Dimension	D	$D + 1/2$	$D + 1$	$D + 3/2$	$D + 2$	$D + 5/2$	$D + 3$	$D + 7/2$	$D + 4$
R – charge	R	$R + 1/2$	$R + 1$	$R + 3/2$	$R + 2$	$R + 5/2$	$R + 3$	$R + 7/2$	$R + 4$

Table 1.1: R -charge content of a supermultiplet

The states at each level can be classified under the Lorentz group and the $SO(4) \sim SU(2) \times SU(2)$ subgroup of the R -symmetry group, which is unbroken after we have fixed the $SO(2)$ R -charge. For instance, the 28 states at level 2 decompose under $SO(4)_{Lor} \times SO(4)_R$ as $(6, 1) + (1, 6) + (4, 4)$. For the present, the most important point is that, given the dimension of one operator at one level, we can infer the dimensions of all other operators in the supermultiplet.

By working in this fashion we can generate complete anomalous dimension spectra of all two-impurity operators. The results obtained in this manner agree with work originally completed by Beisert in [28]. We will summarize these results here, adding some further useful information that emerges from our own $SU(4)$ analysis. The supermultiplet of interest is based on the set of scalars $\sum_A \text{tr}(\phi^A Z^p \phi^A Z^{R-p})$, the operator class we have denoted by $\bar{T}_{R+2}^{(0)}$. According to (1.1.20), the spectrum of $\Delta = D - R$ eigenvalues associated with this operator basis is

$$\Delta(\bar{T}_{R+2}^{(0)}) = 2 + \frac{g_{YM}^2 N_c}{\pi^2} \sin^2 \left(\frac{n\pi}{R+3} \right) \rightarrow 2 + \frac{g_{YM}^2 N_c}{R^2} n^2 \left(1 - \frac{6}{R} + O(R^{-2}) \right). \quad (1.2.1)$$

The remaining scalar operators $\bar{T}_{R+2}^{(\pm)}$ are included in the supermultiplet and the dimension formulas are expressed in terms of the R -charge of the lowest-dimension member. It turns out that (1.2.1) governs *all* the operators at *all* levels in the supermultiplet. The results of this program, carried out on the spacetime scalar operators, are summarized in table 1.2.

L	R	$SU(4)$ Irreps	Operator	$\Delta - 2$	Multiplicity
0	R_0	$(0, R_0, 0)$	$\Sigma_A \text{tr} (\phi^A Z^p \phi^A Z^{R_0-p})$	$\frac{g_Y^2 M^{N_c}}{\pi^2} \sin^2(\frac{n\pi}{(R_0)+3})$	$n = 1, \dots, \frac{R_0+1}{2}$
2	$R_0 + 1$	$(0, R_0, 2) + c.c.$	$\text{tr} (\phi^{[i} Z^p \phi^{j]} Z^{R_0+1-p})$	$\frac{g_Y^2 M^{N_c}}{\pi^2} \sin^2(\frac{n\pi}{(R_0+1)+2})$	$n = 1, \dots, \frac{R_0+1}{2}$
4	$R_0 + 2$	$(2, R_0, 2)$	$\text{tr} (\phi^{(i} Z^p \phi^{j)} Z^{R_0+2-p})$	$\frac{g_Y^2 M^{N_c}}{\pi^2} \sin^2(\frac{n\pi}{(R_0+2)+1})$	$n = 1, \dots, \frac{R_0+1}{2}$
4	$R_0 + 2$	$(0, R_0 + 2, 0) \times 2$	$\text{tr} (\chi^{[\alpha} Z^p \chi^{\beta]} Z^{R_0+1-p})$	$\frac{g_Y^2 M^{N_c}}{\pi^2} \sin^2(\frac{n\pi}{(R_0+2)+1})$	$n = 1, \dots, \frac{R_0+1}{2}$
6	$R_0 + 3$	$(0, R_0 + 2, 2) + c.c.$	$\text{tr} (\chi^{(\alpha} Z^p \chi^{\beta)} Z^{R_0+2-p})$	$\frac{g_Y^2 M^{N_c}}{\pi^2} \sin^2(\frac{n\pi}{(R_0+3)+0})$	$n = 1, \dots, \frac{R_0+1}{2}$
8	$R_0 + 4$	$(0, R_0, 0)$	$\text{tr} (\nabla_\mu Z Z^p \nabla^\mu Z Z^{R_0+2-p})$	$\frac{g_Y^2 M^{N_c}}{\pi^2} \sin^2(\frac{n\pi}{(R_0+4)-1})$	$n = 1, \dots, \frac{R_0+1}{2}$

Table 1.2: Dimensions and multiplicities of spacetime scalar operators

The supermultiplet contains operators that are not spacetime scalars (i.e., that transform nontrivially under the $SU(2, 2)$ conformal group) and group theory determines at what levels in the supermultiplet they must lie. A representative sampling of data on such operators (extracted from Beisert's paper) is collected in table 1.3. We have worked out neither the $SU(4)$ representations to which these lowest- Δ operators belong nor their precise multiplicities. The ellipses indicate that the operators in question contain further monomials involving fermion fields (so that they are not uniquely specified by their bosonic content). This information will be useful in consistency checks to be carried out below.

L	R	Operator	$\Delta - 2$	$\Delta - 2 \rightarrow$
2	$R_0 + 1$	$\text{tr}(\phi^i Z^p \nabla_\mu Z Z^{R_0-p}) + \dots$	$\frac{g_{YM}^2 N_c}{\pi^2} \sin^2\left(\frac{n\pi}{(R_0+1)+2}\right)$	$\frac{g_{YM}^2 N_c}{R_0^2} n^2 \left(1 - \frac{4}{R_0}\right)$
4	$R_0 + 2$	$\text{tr}(\phi^i Z^p \nabla_\mu Z Z^{R_0+1-p})$	$\frac{g_{YM}^2 N_c}{\pi^2} \sin^2\left(\frac{n\pi}{(R_0+2)+1}\right)$	$\frac{g_{YM}^2 N_c}{R_0^2} n^2 \left(1 - \frac{2}{R_0}\right)$
4	$R_0 + 2$	$\text{tr}(\nabla_{(\mu} Z Z^p \nabla_{\nu)} Z Z^{R_0-p})$	$\frac{g_{YM}^2 N_c}{\pi^2} \sin^2\left(\frac{n\pi}{(R_0+2)+1}\right)$	$\frac{g_{YM}^2 N_c}{R_0^2} n^2 \left(1 - \frac{2}{R_0}\right)$
6	$R_0 + 3$	$\text{tr}(\phi^i Z^p \nabla_\mu Z Z^{R_0+2-p}) + \dots$	$\frac{g_{YM}^2 N_c}{\pi^2} \sin^2\left(\frac{n\pi}{R_0+3}\right)$	$\frac{g_{YM}^2 N_c}{R_0^2} n^2 \left(1 - \frac{0}{R_0}\right)$
6	$R_0 + 3$	$\text{tr}(\nabla_{[\mu} Z Z^p \nabla_{\nu]} Z Z^{R_0+1-p})$	$\frac{g_{YM}^2 N_c}{\pi^2} \sin^2\left(\frac{n\pi}{R_0+3}\right)$	$\frac{g_{YM}^2 N_c}{R_0^2} n^2 \left(1 - \frac{0}{R_0}\right)$

Table 1.3: Anomalous dimensions of some operators that are not scalars

Level	0	1	2	3	4	5	6	7	8
Multiplicity	1	8	28	56	70	56	28	8	1
$\delta E \times (R^2/g_{YM}^2 N_c n^2)$	$-6/R$	$-5/R$	$-4/R$	$-3/R$	$-2/R$	$-1/R$	0	$1/R$	$2/R$

Table 1.4: Predicted energy shifts of two-impurity string states

The complete dimension spectrum of operators with R -charge R at level L in the supermultiplet are given by the general formula (valid for large R and fixed n):

$$\begin{aligned}
\Delta_n^{R,L} &= 2 + \frac{g_{YM}^2 N_c}{\pi^2} \sin^2\left(\frac{n\pi}{R+3-L/2}\right) \\
&= 2 + \frac{g_{YM}^2 N_c}{R^2} n^2 \left(1 - \frac{6-L}{R} + O(R^{-2})\right). \quad (1.2.2)
\end{aligned}$$

It should be emphasized that, for fixed R , the operators associated with different levels are actually coming from *different* supermultiplets; this is why they have different dimensions! As mentioned before, we can also precisely identify transformation properties under the Lorentz group and under the rest of the R -symmetry group of the degenerate states at each level. This again leads to useful consistency checks, and we will elaborate on this when we analyze the eigenstates of the string worldsheet Hamiltonian.

Chapter 2

A virial approach to operator dimensions

In the previous chapter we reviewed how the problem of computing operator dimensions in the planar limit of large- N $\mathcal{N} = 4$ SYM theory maps to that of diagonalizing the Hamiltonian of certain quantum mechanical systems. Calculating operator dimensions is equivalent to finding the eigenvalue spectrum of spin chain Hamiltonians, and various established techniques associated with integrable systems (most notably the Bethe ansatz) have proved useful in this context (for a general review of the Bethe ansatz method, see [67]). The utility of this approach was first demonstrated by Minahan and Zarembo in [39]. For operators with two R -charge impurities, the spin chain spectra can be computed exactly via the Bethe ansatz. For three- or higher-impurity operators, however, the Bethe equations have only been solved perturbatively near the limit of infinite chain length [32, 39, 68]. Furthermore, at higher-loop order in λ , the spin chain Hamiltonians typically acquire long-range or non-nearest-neighbor interactions for which a general Bethe ansatz may not be available. For example, while the action of the spin chain Hamiltonian in the “closed $\mathfrak{su}(2|3)$ ” sector is known to three-loop order [33], the corresponding long-range Bethe ansatz is not known (though it may well exist). (See [52] for a more recent approach to deriving Bethe ansatz equations.) A long-range Bethe ansatz does exist for the particularly simple “closed $\mathfrak{su}(2)$ ” sector of the theory [34, 37], and our methods will provide a useful cross-check on these approaches to gauge theory anomalous dimensions at higher

order in the 't Hooft parameter $\lambda = g_{\text{YM}}^2 N_c$.

In this chapter we will present a virial approach to the spin chain systems of $\mathcal{N} = 4$ SYM theory. The generic spin chain Hamiltonian acts on single-impurity pseudoparticles as a lattice Laplacian and higher N -body interactions among pseudoparticles are suppressed relative to the one-body pseudoparticle energy by inverse powers of the lattice length K . Surprisingly, this expansion of the spin chain Hamiltonian is truncated at $O(K^{-3})$ in certain subsectors of the theory, allowing straightforward eigenvalue calculations that are exact in the chain length for operators with more than two R -charge impurities. Furthermore, since the goal is to eventually compare anomalous dimensions with $1/J$ energy corrections to corresponding string states near the pp-wave limit of $AdS_5 \times S^5$, and because the string angular momentum J is proportional to the lattice length K , this virial expansion is precisely what is needed to devise a practical method for testing the AdS/CFT correspondence at any order in the gauge theory loop expansion for an arbitrary number of R -charge (or worldsheet) impurities.

We will focus on three particular closed sectors of the theory, each labeled by the subalgebra of the full superconformal algebra that characterizes the spin variables of the equivalent spin chain system. Specifically, there are two sectors spanned by bosonic operators and labeled by $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ subalgebras plus an $\mathfrak{su}(2|3)$ sector which includes fermionic operators. Section 2.1 is dedicated to an analysis of the bosonic $\mathfrak{su}(2)$ closed sector to three-loop order in λ . In Section 2.2 we analyze an $\mathfrak{su}(1|1)$ subsector of the closed $\mathfrak{su}(2|3)$ system to three-loop order. The spin chain Hamiltonian in the bosonic $\mathfrak{sl}(2)$ sector has previously been determined to one loop, and we analyze this system in Section 2.3.

2.1 The $\mathfrak{su}(2)$ sector

Single-trace operators in the closed $\mathfrak{su}(2)$ sector are constructed from two complex scalar fields of $\mathcal{N} = 4$ SYM, typically denoted by Z and ϕ . Under the $SO(6) \simeq$

$U(1)_R \times SO(4)$ decomposition of the full $SU(4)$ R -symmetry group, the Z fields are charged under the scalar $U(1)_R$ component and ϕ is a particular scalar field carrying zero R -charge. The basis of length- K operators in the planar limit is constructed from single-trace monomials with I impurities and total R -charge equal to $K - I$:

$$\text{tr}(\phi^I Z^{K-I}) , \quad \text{tr}(\phi^{I-1} Z \phi Z^{K-I-1}) , \quad \text{tr}(\phi^{I-2} Z \phi^2 Z^{K-I-1}) , \quad \dots . \quad (2.1.1)$$

The statement that this sector of operators is “closed” means simply that the anomalous dimension operator can be diagonalized on this basis, at least to leading order in large N_c [31, 69].

The heart of the spin chain approach is the proposition that there exists a one-dimensional spin system whose Hamiltonian can be identified with the large- N_c limit of the anomalous dimension operator acting on this closed subspace of operators [39]. Since the anomalous dimensions are perturbative in the 't Hooft coupling λ , it is natural to expand the $\mathfrak{su}(2)$ spin chain Hamiltonian in powers of λ as well:

$$H_{\mathfrak{su}(2)} = I + \sum_n \left(\frac{\lambda}{8\pi^2} \right)^n H_{\mathfrak{su}(2)}^{(2n)} . \quad (2.1.2)$$

Comparison with the gauge theory has shown that successive terms in the expansion of the Hamiltonian have a remarkably simple structure: the one-loop-order Hamiltonian $H_{\mathfrak{su}(2)}^{(2)}$ is built out of permutations of pairs of nearest-neighbor fields and, at n^{th} order, the Hamiltonian permutes among themselves fields that are at most n lattice sites apart. This is a universal structure that leads to remarkable simplifications in the various closed sectors of the theory [32].

Beisert, Kristjansen and Staudacher [31] have introduced the following useful notation for products of permutations acting on operators separated by an arbitrary number of lattice sites:

$$\{n_1, n_2, \dots\} = \sum_{k=1}^K P_{k+n_1, k+n_1+1} P_{k+n_2, k+n_2+1} \dots , \quad (2.1.3)$$

where $P_{i,j}$ simply exchanges fields on the i^{th} and j^{th} lattice sites on the chain. The upshot of the gauge theory analysis is that the equivalent spin chain Hamiltonian for the $\mathfrak{su}(2)$ sector can be written in a rather compact form in terms of this notation. The result, correct to three-loop order, is (see [31] for details)

$$H_{\mathfrak{su}(2)}^{(2)} = 2(\{\} - \{0\}) , \quad (2.1.4)$$

$$H_{\mathfrak{su}(2)}^{(4)} = 2(-4\{\} + 6\{0\} - (\{0, 1\} + \{1, 0\})) , \quad (2.1.5)$$

$$H_{\mathfrak{su}(2)}^{(6)} = 4[15\{\} - 26\{0\} + 6(\{0, 1\} + \{1, 0\}) + \{0, 2\} - (\{0, 1, 2\} + \{2, 1, 0\})] . \quad (2.1.6)$$

(Note that $\{\}$ is just the identity operator.) The form of the three-loop term $H_{\mathfrak{su}(2)}^{(6)}$ was first conjectured in [31] based on integrability restrictions and BMN scaling; this conjecture was later corroborated by direct field-theoretic methods in [33] (see also [30] for relevant discussion on this point). Our goal is to develop practical methods for finding the eigenvalue spectrum of the spin chain Hamiltonian for various interesting cases.

2.1.1 One-loop order

We start at one-loop order with $H_{\mathfrak{su}(2)}^{(2)}$ in eqn. (2.1.4), which provides a natural ‘position-space’ prescription for constructing matrix elements in an I -impurity basis of operators. As an explicit example, we consider first the basis of two-impurity operators of length $K = 8$:

$$\text{tr}(\phi^2 Z^6) , \quad \text{tr}(\phi Z \phi Z^5) , \quad \text{tr}(\phi Z^2 \phi Z^4) , \quad \text{tr}(\phi Z^3 \phi Z^3) . \quad (2.1.7)$$

It is easy to see that the one-loop Hamiltonian mixes the four elements of this basis according to the matrix

$$H_{\mathfrak{su}(2)}^{(2)} = \begin{pmatrix} 2 & -2 & 0 & 0 \\ -2 & 4 & -2 & 0 \\ 0 & -2 & 4 & -2\sqrt{2} \\ 0 & 0 & -2\sqrt{2} & 4 \end{pmatrix}. \quad (2.1.8)$$

This matrix generalizes to arbitrary K and it is simple to show that the two-impurity one-loop eigenvalues of $H_{\mathfrak{su}(2)}^{(2)}$ are given by the formula [28]

$$E_{\mathfrak{su}(2)}^{(2)} = 8 \sin^2 \left(\frac{\pi n}{K-1} \right), \quad n = 0, \dots, n_{\max} = \begin{cases} (K-2)/2, & K \text{ even} \\ (K-3)/2, & K \text{ odd} \end{cases}. \quad (2.1.9)$$

Note that if the denominator $K-1$ were replaced by K , the above expression would agree with the usual lattice Laplacian energy for a lattice of length K . The difference amounts to corrections to the free Laplacian of higher order in $1/K$ and we will seek to understand the physical origin of such corrections in what follows.

To compare gauge theory predictions with $1/J$ corrections to the three-impurity spectrum of the string theory on $AdS_5 \times S^5$, we need to determine the large- K behavior of the three-impurity spin chain spectrum. We are primarily interested in systems with few impurities compared to the length of the spin chain and we expect that impurity interaction terms in the Hamiltonian will be suppressed by powers of the impurity density (i.e., inverse powers of the lattice length). This suggests that we develop a virial expansion of the spin chain Hamiltonian in which the leading-order term in $1/K$ gives the energy of free pseudoparticle states on the lattice (labeled by lattice momentum mode numbers as in the two-impurity spectrum eqn. (2.1.9)) and higher $1/K$ corrections come from N -body interactions described by vertices V_N . A reasonable guess about how the N -body interactions should scale with $1/K$ suggests

that we can write the one-loop-order energy for I impurities in the form

$$E(\{n_i\}) = I + \frac{\lambda}{2\pi^2} \sum_{i=1}^I \sin^2 \frac{n_i \pi}{K} + \sum_{N=2}^{2I} \frac{\lambda}{K^{2N-1}} V_{N\text{-body}}(n_1, \dots, n_I) + \dots, \quad (2.1.10)$$

where the leading-order contribution I measures the naive dimension minus R -charge, the next term is the lattice Laplacian energy of I non-interacting pseudoparticles and the $1/K$ corrections account for interactions between pseudoparticles (which may depend on the lattice momenta mode numbers n_i). In the many-body approach, one would try to derive such energy expressions by rewriting the Hamiltonian in terms of creation/annihilation operators $b_{n_i}, b_{n_i}^\dagger$ for the pseudoparticles (commuting or anticommuting as appropriate). The N -body interaction vertex would generically be written in terms of the b, b^\dagger as

$$V_N = \sum_{n_i, m_i} \delta_{n_1+\dots+n_N, m_1+\dots+m_N} f_N(\{n_i\}, \{m_i\}) \prod_{i=1}^N b_{n_i}^\dagger \prod_{i=1}^N b_{m_i}, \quad (2.1.11)$$

where $f_N(\{n_i\}, \{m_i\})$ is some function of the lattice momenta and the Kronecker delta enforces lattice momentum conservation. One has to determine the functions f_N by matching the many-body form of the Hamiltonian to exact spin chain expressions such as eqn. (2.1.4). We will see that, once the Hamiltonian is in many-body form, it is straightforward to obtain a density expansion of the higher-impurity energy eigenvalues.

The discussion so far has been in the context of one-loop gauge theory physics, but the logic of the virial expansion should be applicable to the general case. To include higher-loop order physics we must do two things: a) generalize the functions $f_N(\{n_i\}, \{m_i\})$ defining the multi-particle interaction vertices to power series in λ and b) allow the free pseudoparticle kinetic energies themselves to become power series in λ . We will be able to carry out the detailed construction of the higher-loop virial Hamiltonian in a few well-chosen cases. To match this expansion at n -loop order in λ to the corresponding loop order (in the modified 't Hooft coupling $\lambda' = g_{\text{YM}}^2 N_c / J^2$)

in the string theory, we need to determine the Hamiltonian to $O(K^{-(2n+1)})$ in this virial expansion. (The first curvature correction to the pp-wave string theory at one loop, for example, appears at $O(\lambda/J)$ or, in terms of gauge theory parameters, at $O(\lambda/K^3)$.) Auspiciously, it will turn out that this virial expansion in the $\mathfrak{su}(2)$ sector is truncated at small orders in $1/K$, allowing for simple eigenvalue calculations that are exact in K (although perturbative in λ).

The first step toward obtaining the desired virial expansion is to recast the spin chain Hamiltonian $H_{\mathfrak{su}(2)}$, which is initially expressed in terms of permutation operators, in terms of a creation and annihilation operator algebra. We begin by introducing the spin operators

$$S^\pm = \frac{1}{2}(\sigma_x \pm i\sigma_y) \ , \quad S^z = \frac{1}{2}\sigma_z \ , \quad (2.1.12)$$

where $\vec{\sigma}$ are the Pauli matrices and S_j^\pm , S_j^z act on a two-dimensional spinor space at the j^{th} lattice site in the chain. In this setting the Z and ϕ fields are understood to be modeled by up and down spins on the lattice. The nearest-neighbor permutation operator $P_{i,i+1}$ can be written in terms of spin operators as

$$P_{i,i+1} = S_i^+ S_{i+1}^- + S_i^- S_{i+1}^+ + 2S_i^z S_{i+1}^z + \frac{1}{2} \ , \quad (2.1.13)$$

and the one-loop Hamiltonian in eqn. (2.1.4) can be written as

$$H_{\mathfrak{su}(2)}^{(2)} = - \sum_{j=1}^K (S_j^+ S_{j+1}^- + S_j^- S_{j+1}^+) - 2 \sum_{j=1}^K S_j^z S_{j+1}^z + \frac{1}{2} \ . \quad (2.1.14)$$

A Jordan-Wigner transformation can now be used to express the spin generators in terms of anti-commuting creation and annihilation operators (anticommuting because each site can be either unoccupied (Z) or occupied once (ϕ)). A pedagogical introduction to this technique can be found in [70]. The explicit transformation is

$$S_j^+ = b_j^\dagger K(j) = K(j) b_j^\dagger \ ,$$

$$\begin{aligned}
S_j^- &= K(j)b_j = b_jK(j) , \\
S_j^z &= b_j^\dagger b_j - 1/2 ,
\end{aligned}
\tag{2.1.15}$$

where the Klein factors

$$K(j) = \exp \left(i\pi \sum_{k=1}^{j-1} b_k^\dagger b_k \right) \tag{2.1.16}$$

serve to ensure that spin operators on different sites commute, despite the anticommuting nature of the b_j . The functions $K(j)$ are real, Abelian and, for $j \leq k$,

$$[K(j), \mathbf{S}_k] = 0 . \tag{2.1.17}$$

The operators b_j^\dagger and b_j can therefore be written as

$$b_j^\dagger = S_j^+ K(j) , \quad b_j = S_j^- K(j) , \tag{2.1.18}$$

and we easily verify that they satisfy the standard anticommutation relations

$$\{b_j, b_k^\dagger\} = \delta_{jk} , \quad \{b_j^\dagger, b_k^\dagger\} = \{b_j, b_k\} = 0 . \tag{2.1.19}$$

Cyclicity on the lattice requires that $\mathbf{S}_{K+1} = \mathbf{S}_1$, a condition that can be enforced by the following boundary condition on the creation and annihilation operators:

$$b_{K+1} = (-1)^{I+1} b_1 , \quad I \equiv \sum_{j=1}^K b_j^\dagger b_j , \tag{2.1.20}$$

where the integer I counts the number of spin chain impurities. In this chapter we will be primarily interested in analyzing spin chains with three impurities. The two-impurity problem can usually be solved more directly and, although the techniques presented here are certainly applicable, going to four impurities introduces unnecessary complications. We will henceforth impose the boundary conditions in

eqn. (2.1.20) for odd impurity number only. We can use all of this to re-express eqn. (2.1.14) in creation and annihilation operator language, with the result

$$H_{\mathfrak{su}(2)}^{(2)} = \sum_{j=1}^K \left(b_j^\dagger b_j + b_{j+1}^\dagger b_{j+1} - b_{j+1}^\dagger b_j - b_j^\dagger b_{j+1} + 2 b_j^\dagger b_{j+1}^\dagger b_j b_{j+1} \right) . \quad (2.1.21)$$

Converting to momentum space via the usual Fourier transform

$$b_j = \frac{1}{\sqrt{K}} \sum_{p=0}^{K-1} e^{-\frac{2\pi i j}{K} p} \tilde{b}_p \quad (2.1.22)$$

yields

$$H_{\mathfrak{su}(2)}^{(2)} = 4 \sum_{p=0}^{K-1} \sin^2 \left(\frac{\pi p}{K} \right) \tilde{b}_p^\dagger \tilde{b}_p + \frac{2}{K} \sum_{p,q,r,s=0}^{K-1} e^{\frac{2\pi i (q-s)}{K}} \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r \tilde{b}_s \delta_{p+q,r+s} . \quad (2.1.23)$$

This is a rather standard many-body Hamiltonian: it acts on a Fock space of momentum eigenstate pseudoparticles, contains a one-body pseudoparticle kinetic energy term and a two-body pseudoparticle interaction (the latter having the critical property that it conserves the number of pseudoparticles). Note that the Hamiltonian terminates at two-body interactions, a fact that will simplify the virial expansion of the energy spectrum. This termination is a consequence of the fact that the one-loop Hamiltonian contains only nearest-neighbor interactions and that lattice sites can only be once-occupied.

Because the pseudoparticle (or impurity) number is conserved by the interaction, three-impurity eigenstates of the Hamiltonian must lie in the space spanned by

$$\tilde{b}_{k_1}^\dagger \tilde{b}_{k_2}^\dagger \tilde{b}_{k_3}^\dagger |K\rangle , \quad k_1 + k_2 + k_3 = 0 \pmod{K} , \quad (2.1.24)$$

where the ground state $|K\rangle$ is identified with the zero-impurity operator $\text{tr}(Z^K)$ and the condition of vanishing net lattice momentum arises from translation invariance on the spin chain (which in turn arises from the cyclicity of the single-trace operators

in the operator basis). As a concrete example, the basis of three-impurity states of the $K = 6$ $\mathfrak{su}(2)$ spin chain is

$$\tilde{b}_0^\dagger \tilde{b}_1^\dagger \tilde{b}_5^\dagger |K\rangle, \quad \tilde{b}_0^\dagger \tilde{b}_2^\dagger \tilde{b}_4^\dagger |K\rangle, \quad \tilde{b}_1^\dagger \tilde{b}_2^\dagger \tilde{b}_3^\dagger |K\rangle, \quad \tilde{b}_3^\dagger \tilde{b}_4^\dagger \tilde{b}_5^\dagger |K\rangle, \quad (2.1.25)$$

and the matrix elements of the Hamiltonian (2.1.23) in this basis are easily computed:

$$H_{\mathfrak{su}(2)}^{(2)} = \begin{pmatrix} \frac{1}{3} & -1 & \frac{1}{3} & \frac{1}{3} \\ -1 & 3 & -1 & -1 \\ \frac{1}{3} & -1 & \frac{19}{3} & \frac{1}{3} \\ \frac{1}{3} & -1 & \frac{1}{3} & \frac{19}{3} \end{pmatrix}. \quad (2.1.26)$$

The first-order perturbation theory corrections to the three-impurity operator anomalous dimensions are the eigenvalues of this matrix.

The construction and diagonalization of the Hamiltonian matrix on the degenerate basis of three-impurity operators can easily be carried out for larger K . The results of doing this¹ for lattice sizes out to $K = 40$ are displayed in figure 2.1. According to

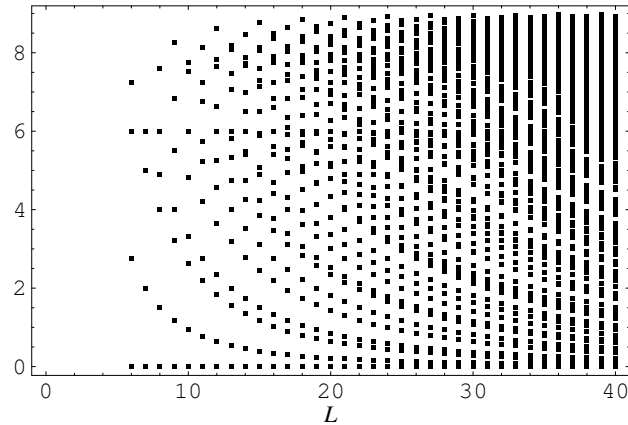


Figure 2.1: One-loop $\mathfrak{su}(2)$ spin chain spectrum vs. lattice length K ($6 \leq K \leq 40$)

¹Using the position- or momentum-space formalism is purely a matter of convenience. In practice we have found that for all sectors the momentum-space treatment is computationally much more efficient. The large- K extrapolations of both methods can be checked against each other, and we of course find that they are in agreement.

eqn. (2.1.10), we expect the eigenvalues of $H_{\mathfrak{su}(2)}^{(2)}$ to scale for large K according to

$$E_K(\{k_i\}) = \frac{\lambda}{K^2} E_{\mathfrak{su}(2)}^{(1,2)}(\{k_i\}) + \frac{\lambda}{K^3} E_{\mathfrak{su}(2)}^{(1,3)}(\{k_i\}) + O(\lambda K^{-4}) . \quad (2.1.27)$$

The scaling coefficients $E_{\mathfrak{su}(2)}^{(1,2)}$ and $E_{\mathfrak{su}(2)}^{(1,3)}$ can easily be extracted from the data displayed in figure 2.1 by fitting the spectral curves to large-order polynomials in $1/K$ (a similar treatment was used in [29]). The results of this procedure are recorded for several low-lying levels in the spectrum (excluding zero eigenvalues) in table 2.1. Subtracting the small errors, we claim that we have the following simple predictions

$E_{\mathfrak{su}(2)}^{(1,2)}$	$E_{\mathfrak{su}(2)}^{(1,3)}$	$E_{\mathfrak{su}(2)}^{(1,3)}/E_{\mathfrak{su}(2)}^{(1,2)}$	Lattice Momenta (k_1, k_2, k_3)
$1 + 2.6 \times 10^{-9}$	$2 - 4.9 \times 10^{-7}$	$2 - 5.0 \times 10^{-7}$	$(1, 0, -1)$
$3 + 4.6 \times 10^{-9}$	$7 - 8.8 \times 10^{-7}$	$7/3 - 3.0 \times 10^{-7}$	$(1, 1, -2)$
$3 + 4.6 \times 10^{-9}$	$7 - 8.8 \times 10^{-7}$	$7/3 - 3.0 \times 10^{-7}$	$(-1, -1, 2)$
$4 + 6.0 \times 10^{-9}$	$8 - 1.1 \times 10^{-6}$	$2 - 2.9 \times 10^{-7}$	$(2, 0, -2)$
$7 + 3.2 \times 10^{-8}$	$14 - 7.1 \times 10^{-6}$	$2 - 1.0 \times 10^{-6}$	$(1, 2, -3)$
$7 + 3.2 \times 10^{-8}$	$14 - 7.1 \times 10^{-6}$	$2 - 1.0 \times 10^{-6}$	$(-1, -2, 3)$
$9 + 2.2 \times 10^{-7}$	$18 - 5.1 \times 10^{-5}$	$2 - 5.7 \times 10^{-6}$	$(3, 0, -3)$
$12 + 5.7 \times 10^{-5}$	$28 + 3.8 \times 10^{-3}$	$7/3 - 1.4 \times 10^{-3}$	$(2, 2, -4)$
$12 + 5.7 \times 10^{-5}$	$28 + 3.8 \times 10^{-3}$	$7/3 - 1.4 \times 10^{-3}$	$(-2, -2, 4)$
$13 - 5.6 \times 10^{-5}$	$26 - 3.8 \times 10^{-3}$	$2 + 1.3 \times 10^{-3}$	$(1, 3, -4)$
$13 - 5.6 \times 10^{-5}$	$26 - 3.8 \times 10^{-3}$	$2 + 1.3 \times 10^{-3}$	$(-1, -3, 4)$

Table 2.1: Scaling limit of three-impurity $\mathfrak{su}(2)$ numerical spectrum at one loop in λ

for the large- K $\mathfrak{su}(2)$ expansion coefficients $E_{\mathfrak{su}(2)}^{(1,3)}$ and $E_{\mathfrak{su}(2)}^{(1,2)}$:

$$\begin{aligned}
E_{\mathfrak{su}(2)}^{(1,2)} &= (k_1^2 + k_2^2 + k_3^2)/2 , & k_1 + k_2 + k_3 &= 0 , \\
E_{\mathfrak{su}(2)}^{(1,3)}/E_{\mathfrak{su}(2)}^{(1,2)} &= 2 , & (k_1 \neq k_2 \neq k_3) , \\
E_{\mathfrak{su}(2)}^{(1,3)}/E_{\mathfrak{su}(2)}^{(1,2)} &= \frac{7}{3} , & (k_1 = k_2, k_3 = -2k_1) .
\end{aligned} \quad (2.1.28)$$

Note the slight annoyance that we must distinguish the case where all mode indices

are unequal from the case where two indices are equal and different from the third. The last column of table 2.1 displays the choice of indices $\{k_i\}$ that best fit each spectral series. As the lattice momenta increase, higher-order $1/K$ corrections to the spectrum become stronger and more data will be required to maintain a given level of precision of the polynomial fit. This effect can be seen directly in the extrapolated eigenvalues in table 2.1.

We also note that the spectrum in table 2.1 exhibits a degeneracy of eigenstates whose momentum labels are related by an overall sign flip (a symmetry that is implemented on the operator basis by a parity operator P that reverses the ordering of all fields within the trace). This degeneracy among “parity pairs” of gauge theory operators was observed in [31], where it was shown that it arises as a consequence of integrability (which can, in turn, be used to constrain the form of the Hamiltonian at higher loop order [30]). See [71] for further discussion on the implications of this degeneracy.

To corroborate these results we turn to the one-loop Bethe ansatz for the Heisenberg spin chain. The Bethe ansatz for chains of spins in arbitrary representations of arbitrary simple Lie groups was developed some time ago [72] (see also [73] for an extension to supersymmetric spin chains) and applied only recently to the specific case of the dilatation operator of $\mathcal{N} = 4$ SYM [32, 39]. In the notation of [32], the Bethe equations are expressed in terms of the so-called Bethe roots (or rapidities) u_i associated with the various impurity insertions in the single-trace ground state $\text{tr}(Z^K)$. In a one-dimensional dynamical interpretation, the impurities are pseudoparticle excitations and the roots parameterize in some fashion the lattice momenta of the pseudoparticles. The index i in the Bethe root u_i runs over the total number I of impurities. A second index $q_i = 1, \dots, 7$ is used to associate each of the I Bethe roots with a particular simple root of the $\mathfrak{sl}(4|4)$ symmetry algebra associated with $\mathcal{N} = 4$ SYM. The Bethe ansatz then takes the form (see [32] and references therein

for further details)

$$\left(\frac{u_i + \frac{i}{2} V_{q_i}}{u_i - \frac{i}{2} V_{q_i}} \right)^K = \prod_{j \neq i}^I \left(\frac{u_i - u_j + \frac{i}{2} M_{q_i, q_j}}{u_i - u_j - \frac{i}{2} M_{q_i, q_j}} \right), \quad (2.1.29)$$

where V_{q_i} denotes the q_i^{th} Dynkin coefficient of the spin representation and M is the Cartan matrix of the algebra. To be slightly more specific, if α_{q_i} are the root vectors associated with the nodes of the Dynkin diagram and μ is the highest weight of the spin representation, then the Dynkin coefficient (for a bosonic algebra) is $V_{q_i} = 2 \alpha^{(q_i)} \cdot \mu / (\alpha^{(q_i)})^2$ and the elements of the Cartan matrix are $M_{q_i, q_j} = 2 \alpha^{(q_i)} \cdot \alpha^{(q_j)} / (\alpha^{(q_j)})^2$ (note that diagonal elements $M_{q_i, q_i} = 2$). (For superalgebras see, e.g., [74, 75].) Furthermore, since the spin chain systems of interest to us are cyclic and carry no net momentum (analogous to the level-matching condition in the string theory), the Bethe roots u_i are subject to the additional constraint

$$1 = \prod_i^I \left(\frac{u_i + \frac{i}{2} V_{q_i}}{u_i - \frac{i}{2} V_{q_i}} \right). \quad (2.1.30)$$

Having found a set of Bethe roots u_i that solve the above equations, the corresponding energy eigenvalue (up to an overall additive constant; see, e.g., [32]) is given by

$$E = \sum_{j=1}^I \left(\frac{V_{q_j}}{u_j^2 + V_{q_j}^2/4} \right). \quad (2.1.31)$$

In the current application all impurities are of the same type (i.e., carry the same Dynkin label), so the index q_i can be ignored. It is worth noting, however, that the Dynkin coefficient V_{q_i} can vanish, in which case the associated Bethe roots do not contribute directly to the energy.

The Bethe equations are typically exactly soluble for the case of two identical impurities (i.e., two Bethe roots u_1, u_2 associated with the same simple root of the algebra). The two-impurity $\mathfrak{su}(2)$ Bethe equations, for example, yield solutions that reproduce the familiar two-impurity anomalous dimension formula noted above in

eqn. (2.1.9) (see [32, 39] for further examples). For three and higher impurities, however, exact solutions are not known. Since we are ultimately interested in comparing with string theory predictions at large values of the S^5 angular momentum J , an alternate approach is to solve the Bethe equations perturbatively in small $1/K$. Experience shows that, in the limit where we can neglect interactions between excitations (or impurities), the Bethe roots are simply the inverse of the conserved momentum carried by the impurities. With a little work, one can show that the Bethe ansatz conditions, eqns. (2.1.29, 2.1.30), can be solved order-by-order in a large- K expansion:

$$u_i = \frac{1}{2\pi k_i} \left(K + A_i \sqrt{K} + B_i + \dots \right) , \quad (2.1.32)$$

where $0 < k_i < K$ is the usual integer lattice momentum. The half-integer powers of K may or may not be present in eqn. (2.1.32): they are needed to deal with special kinematic situations (such as when a pair of impurities has the same lattice momentum) where the integral power expansion would be singular. The eigenvalues of the spin chain (or the anomalous dimensions of the corresponding gauge theory operator) are then obtained as a power series in $1/K$ by substituting the expansion of the Bethe roots into eqn. (2.1.31). This is the approach introduced by Minahan and Zarembo for the $\mathfrak{so}(6)$ spin chain in [39]. Since we wish to carry out similar calculations at higher orders in λ , we will review this methodology at one-loop order for the specific case of three identical impurities in the $\mathfrak{su}(2)$ spin chain. (Since the $\mathfrak{su}(2)$ chain is a subsector of the $\mathfrak{so}(6)$ system studied in [39], the three-impurity Bethe ansatz predictions derived here are of course implied by the all-impurity $\mathfrak{so}(6)$ anomalous dimension formula derived in [39] at one loop.)

We now apply this to the closed $\mathfrak{su}(2)$ sector where the Dynkin diagram has a single node, the Cartan matrix is $M_{\mathfrak{su}(2)} = 2$ and the Dynkin coefficient of the fundamental representation is $V_{\mathfrak{su}(2)} = 1$. Consequently, the Bethe equations (2.1.29, 2.1.30) reduce

to

$$\left(\frac{u_i + i/2}{u_i - i/2}\right)^K = \prod_{j \neq i}^I \left(\frac{u_i - u_j + i}{u_i - u_j - i}\right), \quad (2.1.33)$$

$$1 = \prod_i^I \left(\frac{u_i + i/2}{u_i - i/2}\right). \quad (2.1.34)$$

With three or more pseudoparticle excitations, bound-state solutions can arise that satisfy the second equation (2.1.34). These solutions are characterized as having pseudoparticle states sharing the same lattice momenta (e.g., $k_i = k_j$ for the i^{th} and j^{th} roots). The generic solutions to the Bethe equations can therefore be loosely divided into those that do or do not contain bound states. For three impurities with no bound states present ($k_1 \neq k_2 \neq k_3$), eqn. (2.1.34) states that $k_3 = -k_1 - k_2$. The strategy of [39] can then be used to obtain a systematic expansion of $\mathfrak{su}(2)$ Bethe roots in powers of K^{-1} , with the result

$$\begin{aligned} u_1 &= \frac{K-4}{2\pi k_1} + \frac{3k_1}{\pi(k_1 - k_2)(2k_1 + k_2)} + O(K^{-1}), \\ u_2 &= \frac{(K-4)k_1^2 + (K-4)k_1k_2 - 2(K-1)k_2^2}{2\pi k_2(k_1^2 + k_1k_2 - 2k_2^2)} + O(K^{-1}), \\ u_3 &= -\frac{(K-1)k_1^2 - (8-5K)k_1k_2 + 2(K-1)k_2^2}{2\pi(k_1 + k_2)(2k_1 + k_2)(k_1 + 2k_2)} + O(K^{-1}). \end{aligned} \quad (2.1.35)$$

Substituting these roots into the energy formula in eqn. (2.1.30) gives the following expression for the anomalous dimension of the $\mathfrak{su}(2)$ three-impurity operator at one loop:

$$E_{\mathfrak{su}(2)}^{(2)}(k_1, k_2) = \frac{8\pi^2}{K^3} (k_1^2 + k_1k_2 + k_2^2) (K+2) + O(K^{-4}), \quad (k_1 \neq k_2 \neq k_3). \quad (2.1.36)$$

This is in perfect agreement with the results of eqn. (2.1.28) and the numerical gauge theory results in table 2.1. When a single bound state is present the Bethe roots must be altered. Taking, for example, $k_1 = k_2$, the cyclic constraint in eqn. (2.1.34) sets

$k_3 = -2k_1$, and the Bethe roots are

$$\begin{aligned} u_1 &= \frac{-7 + 3i\sqrt{K} + 3K}{6\pi k_1} + O(K^{-1/2}) , \\ u_2 &= -\frac{7 + 3i\sqrt{K} - 3K}{6\pi k_1} + O(K^{-1/2}) , \\ u_3 &= \frac{4 - 3K}{12\pi k_1} + O(K^{-1/2}) . \end{aligned} \quad (2.1.37)$$

In this case the anomalous dimension is predicted to be

$$\begin{aligned} E_{\mathfrak{su}(2)}^{(2)}(k_1) &= \frac{8\pi^2}{K^3} k_1^2 (3K + 7) + O(K^{-4}) , \\ &\quad (k_1 = k_2, \ k_3 = -2k_1) , \end{aligned} \quad (2.1.38)$$

which is again in agreement with the results of eqn. (2.1.28) and table 2.1 (note that the fractional powers of K^{-1} have obligingly canceled out of the final expression for the energy).

2.1.2 Two- and three-loop order

A similar analysis can be performed on the two-loop $\mathfrak{su}(2)$ spin chain Hamiltonian. As before, we use the Jordan-Wigner transformation restricted to an odd-impurity basis of operators to rewrite the two-loop Hamiltonian (2.1.5) in terms of position-space fermionic oscillators, obtaining a result similar to eqn. (2.1.21):

$$\begin{aligned} H_{\mathfrak{su}(2)}^{(4)} &= \sum_{j=1}^K \left\{ -\frac{1}{2} \left[b_{j+2}^\dagger b_j + b_j^\dagger b_{j+2} - 4 \left(b_{j+1}^\dagger b_j + b_j^\dagger b_{j+1} \right) \right] - 3 b_j^\dagger b_j - 4 b_j^\dagger b_{j+1}^\dagger b_j b_{j+1} \right. \\ &\quad \left. + b_{j+1}^\dagger b_{j+2}^\dagger b_j b_{j+1} + b_j^\dagger b_{j+1}^\dagger b_{j+1} b_{j+2} + b_j^\dagger b_{j+2}^\dagger b_j b_{j+2} \right\} . \end{aligned} \quad (2.1.39)$$

Passing to momentum space, we obtain the two-loop analogue of eqn. (2.1.23):

$$H_{\mathfrak{su}(2)}^{(4)} = -8 \sum_{p=0}^{K-1} \sin^4 \left(\frac{p\pi}{K} \right) \tilde{b}_p^\dagger \tilde{b}_p$$

$$+ \frac{1}{K} \sum_{p,q,r,s=0}^{K-1} \left(e^{\frac{2\pi i(q+r)}{K}} + e^{\frac{-2\pi i(p+s)}{K}} + e^{\frac{4\pi i(q-s)}{K}} - 4 e^{\frac{2\pi i(q-s)}{K}} \right) \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r \tilde{b}_s \delta_{p+q,r+s} . \quad (2.1.40)$$

Although the two-loop Hamiltonian includes “long-range” interactions among non-neighboring lattice sites, the momentum-space Hamiltonian (2.1.40) conveniently terminates at two-body interaction terms. An equally important point is that, for fixed momenta p, q, \dots , the one-body (two-body) operators scale as K^{-4} (K^{-5}) for large K (the corresponding scalings for the one-loop Hamiltonian were K^{-2} (K^{-3})). This special relation between density scaling and power of coupling constant is critical for matching to string theory.

We deal with the problem of finding the eigenvalues of the combined one- and two-loop Hamiltonian via Rayleigh-Schrödinger perturbation theory: at each value of the lattice length K we treat the one-loop operator $H_{\mathfrak{su}(2)}^{(2)}$ as a zeroth-order Hamiltonian and regard $H_{\mathfrak{su}(2)}^{(4)}$ as a first-order perturbation. The $O(\lambda^2)$ corrections to the spectrum of $H_{\mathfrak{su}(2)}^{(2)}$ are then found by taking expectation values of the perturbation $H_{\mathfrak{su}(2)}^{(4)}$ in the (numerically-determined) eigenvectors of $H_{\mathfrak{su}(2)}^{(2)}$. This is the recipe for non-degenerate first-order perturbation theory and we might worry that the previously-noted parity-pair degeneracy of the eigenvalues of $H_{\mathfrak{su}(2)}^{(2)}$ would force us to use the rules of degenerate perturbation theory. As discussed in [31, 38, 71], however, parity degeneracy can be traced to the existence of a higher Abelian charge that is conserved to at least three-loop order. This charge can be used to show that the formulas of non-degenerate perturbation theory can be used without modification. The basic observation is that conservation of the Abelian charge guarantees that the matrix element of $H_{\mathfrak{su}(2)}^{(4)}$ between two degenerate eigenstates of $H_{\mathfrak{su}(2)}^{(2)}$ with different eigenvalues of the higher Abelian charge vanishes: this eliminates the vanishing energy-denominator singularities that would otherwise invalidate the non-degenerate first-order perturbation theory formulas (and similar arguments apply to the higher-order cases).

Using this method, we have evaluated the $O(\lambda^2)$ corrections to the spectrum of

anomalous dimensions for lattice sizes from $K = 6$ to $K = 40$. As before, we fit the spectral data to a power series in $1/K$ to read off the leading scaling coefficients of the low-lying eigenvalues. As mentioned in the discussion of the two-loop Hamiltonian (2.1.40), we expect the two-loop eigenvalues to have the following scaling behavior in $1/K$:

$$E_K^{(2)}(\{k_i\}) = \frac{\lambda^2}{K^4} E^{(2,4)}(\{k_i\}) + \frac{\lambda^2}{K^5} E^{(2,5)}(\{k_i\}) + O(\lambda^2 K^{-6}) . \quad (2.1.41)$$

The numerical data confirm that the eigenvalues scale at least as fast as K^{-4} . The resulting numerical values for the leading scaling coefficients of low-lying eigenvalues, $E_{\mathfrak{su}(2)}^{(2,4)}$ and $E_{\mathfrak{su}(2)}^{(2,5)}$, are presented in table 2.2. We thus have the following simple

$E_{\mathfrak{su}(2)}^{(2,4)}$	$E_{\mathfrak{su}(2)}^{(2,5)}$	$E_{\mathfrak{su}(2)}^{(2,5)}/E_{\mathfrak{su}(2)}^{(2,4)}$	(k_1, k_2, k_3)
$-0.25 - 4.6 \times 10^{-9}$	$-2 + 8.0 \times 10^{-7}$	$8 - 3.4 \times 10^{-6}$	$(1, 0, -1)$
$-2.25 - 1.4 \times 10^{-6}$	$-19 + 2.6 \times 10^{-4}$	$76/9 + 1.2 \times 10^{-4}$	$(1, 1, -2)$
$-2.25 - 1.4 \times 10^{-6}$	$-19 + 2.6 \times 10^{-4}$	$76/9 + 1.2 \times 10^{-4}$	$(-1, -1, 2)$
$-4 + 8.3 \times 10^{-7}$	$-32 - 1.1 \times 10^{-4}$	$8 + 3.0 \times 10^{-5}$	$(2, 0, -2)$
$-12.25 - 9.9 \times 10^{-6}$	$-98 + 2.3 \times 10^{-3}$	$8 - 2.0 \times 10^{-4}$	$(1, 2, -3)$
$-12.25 - 9.9 \times 10^{-6}$	$-98 + 2.3 \times 10^{-3}$	$8 - 2.0 \times 10^{-4}$	$(-1, -2, 3)$
$-20.25 + 3.2 \times 10^{-3}$	-161.4	7.97	$(3, 0, -3)$
$-36 - 2.8 \times 10^{-3}$	-304.6	8.46	$(2, 2, -4)$
$-36 - 2.8 \times 10^{-3}$	-304.6	8.46	$(-2, -2, 4)$
$-42.25 + 4.9 \times 10^{-3}$	-337.0	7.97	$(1, 3, -4)$
$-42.25 + 4.9 \times 10^{-3}$	-337.0	7.97	$(-1, -3, 4)$

Table 2.2: Scaling limit of three-impurity $\mathfrak{su}(2)$ numerical spectrum at two loops in λ predictions for the two-loop large- K expansion coefficients:

$$\begin{aligned}
E_{\mathfrak{su}(2)}^{(2,4)} &= -(k_1^2 + k_2^2 + k_3^2)^2/16 , & k_1 + k_2 + k_3 &= 0 , \\
E_{\mathfrak{su}(2)}^{(2,5)}/E_{\mathfrak{su}(2)}^{(2,3)} &= 8 , & (k_1 \neq k_2 \neq k_3) , \\
E_{\mathfrak{su}(2)}^{(2,5)}/E_{\mathfrak{su}(2)}^{(2,3)} &= \frac{76}{9} , & (k_1 = k_2, k_3 = -2k_1) .
\end{aligned} \quad (2.1.42)$$

Once again, the decline in precision as one goes to higher energies is expected. As a consistency check we note that this time we have no freedom to choose the momenta (k_1, k_2, k_3) associated with each state: they have been fixed in the one-loop matching exercise.

The three-loop $\mathfrak{su}(2)$ Hamiltonian (2.1.6) can be dealt with in a similar fashion. The position space operator version of this Hamiltonian is too long to record here, but its momentum space version is fairly compact:

$$\begin{aligned}
H_{\mathfrak{su}(2)}^{(6)} = & 32 \sum_{p=0}^{K-1} \sin^6 \left(\frac{p\pi}{K} \right) \tilde{b}_p^\dagger \tilde{b}_p + \frac{1}{2K} \sum_{p,q,r,s=0}^{K-1} \left\{ -10 e^{\frac{2\pi i(q+r)}{K}} + e^{\frac{2\pi i(2q+r)}{K}} + e^{\frac{2\pi i(q+2r)}{K}} \right. \\
& + e^{\frac{2\pi i(q-3s)}{K}} + e^{\frac{2\pi i(2q-2r-3s)}{K}} + e^{\frac{2\pi i(3q-2r-3s)}{K}} + e^{\frac{2\pi i(q-r-3s)}{K}} + e^{\frac{2\pi i(2q-r-3s)}{K}} \\
& - e^{\frac{2\pi i(q-2s)}{K}} - 10 e^{\frac{2\pi i(q-r-2s)}{K}} - e^{\frac{2\pi i(2q-r-2s)}{K}} - e^{\frac{2\pi i(3q-r-2s)}{K}} - e^{\frac{2\pi i(q+r-2s)}{K}} \\
& + 29 e^{\frac{2\pi i(q-s)}{K}} - 10 e^{\frac{4\pi i(q-s)}{K}} + e^{\frac{6\pi i(q-s)}{K}} - e^{\frac{2\pi i(2q-s)}{K}} + e^{\frac{2\pi i(3q-s)}{K}} \\
& \left. - e^{\frac{2\pi i(q+r-s)}{K}} + e^{\frac{2\pi i(2q+r-s)}{K}} + e^{\frac{2\pi i(q+2r-s)}{K}} \right\} \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r^\dagger \tilde{b}_s \delta_{p+q,r+s} \\
& + \frac{1}{K^2} \sum_{p,q,r,s,t,u=0}^{K-1} \left\{ e^{\frac{2\pi i(q+3r-2t-3u)}{K}} + e^{\frac{2\pi i(q+2r-s-2t-3u)}{K}} \right. \\
& \left. + e^{\frac{2\pi i(2q+3r-t-3u)}{K}} + e^{\frac{2\pi i(q+2r+s-u)}{K}} \right\} \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r^\dagger \tilde{b}_s \tilde{b}_t \tilde{b}_u \delta_{p+q+r,s+t+u} . \quad (2.1.43)
\end{aligned}$$

It contains at most three-body operators and a careful examination of terms shows that, for fixed momenta, the one-body operators scale as K^{-6} , the two-body operators as K^{-7} and so on. We therefore expect the leading scaling coefficients in the $O(\lambda^3)$ eigenvalues to be $E_{\mathfrak{su}(2)}^{(3,6)}$ and $E_{\mathfrak{su}(2)}^{(3,7)}$, to use a by-now-familiar notation. To find the eigenvalues to this order, we continue with the Rayleigh-Schrödinger perturbation theory strategy: the $O(\lambda^3)$ correction to any eigenvalue is the sum of the matrix element of $H_{\mathfrak{su}(2)}^{(6)}$ in the appropriate eigenvector of $H_{\mathfrak{su}(2)}^{(2)}$ plus the second-order sum-over-states contribution of $H_{\mathfrak{su}(2)}^{(4)}$. These two pieces can easily be computed numerically from the explicit Hamiltonian operators at a fixed K . Parity degeneracy and conservation of the higher Abelian charge mentioned above continue to hold, and we can again use non-degenerate perturbation theory formulas to compute the

eigenvalue corrections. We have generated numerical eigenvalue data for lattices from $K = 6$ to $K = 40$ and the large- K scaling coefficients of the low-lying states extracted from those data are given in table 2.3.

$E_{\mathfrak{su}(2)}^{(3,4)}$	$E_{\mathfrak{su}(2)}^{(3,7)}$	$E_{\mathfrak{su}(2)}^{(3,7)}/E_{\mathfrak{su}(2)}^{(3,6)}$	(k_1, k_2, k_3)
0.1250	2.0003	16.003	$(1, 0, -1)$
4.125	58.03	14.07	$(1, 1, -2)$
4.125	58.03	14.07	$(-1, -1, 2)$
7.999	128.2	16.03	$(2, 0, -2)$
49.62	713.3	14.37	$(1, 2, -3)$
49.62	713.3	14.37	$(-1, -2, 3)$
91.15	1, 454	15.96	$(3, 0, -3)$
263.8	3, 739	14.17	$(2, 2, -4)$
263.8	3, 739	14.17	$(-2, -2, 4)$

Table 2.3: Scaling limit of three-impurity $\mathfrak{su}(2)$ numerical spectrum at three loops in λ

A modified Bethe ansatz for the $\mathfrak{su}(2)$ sector of the gauge theory, possibly incorporating all orders of higher-loop physics, has been proposed in [34, 37].² It is an instructive exercise and a useful consistency check on this bold proposal to verify that it reproduces the higher-loop scaling coefficients for three-impurity anomalous dimensions that we have just computed by virial methods (and displayed in tables 2.1, 2.2 and 2.3). For completeness, we briefly summarize the new ansatz, referring the reader to [34] for a detailed account. In the new ansatz, the momenta p_i of the excitations (closely related to the Bethe roots) become functions of λ (as well as K and mode numbers) and are determined by a modified version of eqns. (2.1.33, 2.1.34):

$$e^{iKp_i} = \prod_{j \neq i}^I \frac{\varphi(p_i) - \varphi(p_j) + i}{\varphi(p_i) - \varphi(p_j) - i}, \quad \sum_{i=1}^I p_i = 0. \quad (2.1.44)$$

²The long-range ansatz based on the Inozemtsev spin chain in [37] suffers from improper BMN scaling at four-loop order, a problem that is surmounted in [34]. For further insights into the importance of BMN scaling, see [58].

Dependence on λ enters through the phase function $\varphi(p_i)$, which is defined in terms of the excitation momenta p_i as follows:

$$\varphi(p_i) \equiv \frac{1}{2} \cot(p_i/2) \sqrt{1 + \frac{\lambda}{\pi^2} \sin^2(p_i/2)} . \quad (2.1.45)$$

The energy eigenvalue corresponding to a particular root of these equations is given in terms of the excitation momenta p_i by the formula

$$E_{\mathfrak{su}(2)} = \sum_{i=1}^I \frac{8\pi^2}{\lambda} \left(\sqrt{1 + \frac{\lambda}{\pi^2} \sin^2(p_i/2)} - 1 \right) . \quad (2.1.46)$$

Finding exact solutions of these equations is even more difficult than before, but we can follow the previous strategy of developing an expansion in powers of $1/K$ about non-interacting impurities on an infinite lattice. This is achieved by expanding the excitation momenta p_i according to

$$p_i = \frac{2\pi k_i}{K} + \sum_{n=1} \frac{p_i^{(n)}}{K^{\frac{n+2}{2}}} , \quad (2.1.47)$$

where the integers k_i (subject to the cyclicity constraint $\sum_i k_i = 0$) characterize the non-interacting state about which the expansion is developed. The appearance of half-integer powers of K^{-1} in this expansion is needed to accommodate bound-state solutions to the Bethe equations that arise when some of the momenta k_i are equal. Solutions to the Bethe equation (2.1.44) will determine the expansion coefficients $p_i^{(n)}$ in terms of the mode numbers k_i and ultimately lead to expansions of the energies as power series in K^{-1} , with coefficients that are functions of λ/K^2 .

Explicit results for the K^{-1} expansion of gauge theory operators of arbitrary impurity number, derived by the above method, were presented in [44].³ As usual, expressions are different depending on whether all momenta are unequal or some

³It is important to note that the focus of this chapter is a *different* Bethe ansatz, designed to match the spectrum of the string theory: the gauge theory Bethe ansatz results are derived for comparison purposes.

subset of them is equal. For all mode numbers k_i unequal the I -impurity energy formula in [44] is

$$E_{\text{su}(2)} = K - I + \sum_{i=1}^I \left(\sqrt{1 + \lambda' k_i^2} - \frac{\lambda'}{K - I} \frac{I k_i^2}{\sqrt{1 + \lambda' k_i^2}} \right) - \frac{\lambda'}{K - I} \sum_{\substack{i,j=1 \\ i \neq j}}^I \frac{2k_i^2 k_j}{k_i^2 - k_j^2} \left(k_j + k_i \sqrt{\frac{1 + \lambda' k_j^2}{1 + \lambda' k_i^2}} \right) + O(K^{-2}) , \quad (2.1.48)$$

where we have used $\lambda' = \lambda/J^2 = \lambda/(K - I)^2$ for convenience ($J = K - I$ is the total R -charge). To compare with our virial results, we must further expand in λ ; expanding to first and second order yields the following scaling coefficients (valid for all k_i unequal):

$$\begin{aligned} E_{\text{su}(2)}^{(1,2)} &= k_1^2 + k_1 k_2 + k_2^2 , & E_{\text{su}(2)}^{(1,3)} &= 2(k_1^2 + k_1 k_2 + k_2^2) , \\ E_{\text{su}(2)}^{(2,4)} &= -\frac{1}{4}(q^2 + qr + r^2)^2 , & E_{\text{su}(2)}^{(2,5)} &= -2(q^2 + qr + r^2)^2 . \end{aligned} \quad (2.1.49)$$

These one- and two-loop coefficients match the numerical results presented in tables 2.1 and 2.2 and the analytic string formulas in eqns. (2.1.28, 2.1.42). It is harder to write down a general formula for the many cases in which subsets of momenta are equal but the solution for the particular case of three impurities with a two-excitation bound state ($k_1 = k_2 = n$, $k_3 = -2n$) was also presented in [44]:

$$E_{\text{su}(2)} = K - 3 + 2\sqrt{1 + \lambda' n^2} + \sqrt{1 + \lambda' 4n^2} - \frac{\lambda' n^2}{K - 3} \left(\frac{1}{1 + \lambda' n^2} + \frac{6}{\sqrt{1 + \lambda' n^2}} + \frac{12}{\sqrt{1 + \lambda' 4n^2}} - \frac{8}{\sqrt{1 + \lambda' n^2} \sqrt{1 + \lambda' 4n^2}} \right) . \quad (2.1.50)$$

To compare with the virial results, one must again expand the energy in powers of λ .

Doing so yields the following one- and two-loop bound-state scaling coefficients:

$$\begin{aligned} E_{\mathfrak{su}(2)}^{(1,2)} &= 3 n^2 , & E_{\mathfrak{su}(2)}^{(1,3)} &= 7 n^2 , \\ E_{\mathfrak{su}(2)}^{(2,4)} &= -\frac{9}{4} n^4 , & E_{\mathfrak{su}(2)}^{(2,5)} &= -19 n^4 . \end{aligned} \quad (2.1.51)$$

We easily verify that this agrees with numerical virial results to two-loop order.

The three-loop coefficients obtained by expanding the energy formulas in eqns. (2.1.48, 2.1.50) are given by

$$\begin{aligned} E_{\mathfrak{su}(2)}^{(3,6)} &= \frac{1}{16} \left(2 k_1^6 + 6 k_1^5 k_2 + 15 k_1^4 k_2^2 + 20 k_1^3 k_2^3 \right. \\ &\quad \left. + 15 k_1^2 k_2^4 + 6 k_1 k_2^5 + 2 k_2^6 \right) , \\ E_{\mathfrak{su}(2)}^{(3,7)} &= \frac{1}{4} \left(8 k_1^6 + 24 k_1^5 k_2 + 51 k_1^4 k_2^2 + 62 k_1^3 k_2^3 \right. \\ &\quad \left. + 51 k_1^2 k_2^4 + 24 k_1 k_2^5 + 8 k_2^6 \right) , \end{aligned} \quad (2.1.52)$$

for $(k_1 \neq k_2 \neq k_3)$, and

$$E_{\mathfrak{su}(2)}^{(3,6)} = \frac{33}{8} n^6 , \quad E_{\mathfrak{su}(2)}^{(3,7)} = 58 n^6 , \quad (2.1.53)$$

for the bound-state solution with $(k_1 = k_2 = n, k_3 = -2n)$. The numerical values of these $O(\lambda^3)$ coefficients are tabulated for several low-lying states in the spectrum in table 2.4. The correspondence with table 2.3, which displays the three-loop expansion coefficients extracted from numerical diagonalization of the three-loop Hamiltonian, is good. At this order in the loop expansion higher-order $1/K$ corrections to the spectrum are more important (compared to the one- and two-loop cases), and the numerical extrapolation is less reliable (especially as the lattice momenta increase). The precision can always be improved by including data from larger lattices in the extrapolation. We emphasize that this discussion concerns the different methods of calculation of operator dimensions in the $\mathfrak{su}(2)$ sector only. It seems to us to give useful further evidence that the long-range Bethe ansatz for the $\mathfrak{su}(2)$ sector of the

gauge theory [34] is exact.

$E_{\mathfrak{su}(2)}^{(3,6)}$	$E_{\mathfrak{su}(2)}^{(3,7)}$	$E_{\mathfrak{su}(2)}^{(3,7)}/E_{\mathfrak{su}(2)}^{(3,6)}$	(k_1, k_2, k_3)
0.125	2	16	(1, 0, -1)
4.125	58	14.06	(1, 1, -2)
4.125	58	14.06	(-1, -1, 2)
8	128	16	(2, 0, -2)
49.625	713	14.37	(1, 2, -3)
49.625	713	14.37	(-1, -2, 3)
91.125	1, 458	16	(3, 0, -3)
264	3, 712	14.06	(2, 2, -4)
264	3, 712	14.06	(-2, -2, 4)

Table 2.4: Three-impurity $\mathfrak{su}(2)$ spectrum from the long-range Bethe ansatz at three loops

2.2 A closed $\mathfrak{su}(1|1)$ subsector of $\mathfrak{su}(2|3)$

The three-impurity string theory analysis of [38] identified a fermionic sector of the theory that is diagonalized by string states composed of fermionic excitations projected onto particular four-dimensional subspaces (which transform in an $SU(2)^2 \times SU(2)^2$ notation as a $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ of $SO(4) \times SO(4)$) and symmetrized in their $SO(4) \times SO(4)$ indices. It was also shown that this three-impurity subsector of the theory decouples at all orders in λ .

On the gauge theory side this subsector corresponds to an $\mathfrak{su}(1|1)$ subgroup of the closed $\mathfrak{su}(2|3)$ sector studied by Beisert in [33, 69]. (Supersymmetric integrable $\mathfrak{su}(n|m)$ spin chains have previously been studied in certain condensed-matter applications; see, e.g., [76].) In the present setting the fields of $\mathfrak{su}(2|3)$ consist of three complex scalars ϕ_a and two complex fermions ψ_α . In the closed $\mathfrak{su}(1|1)$ subspace we are restricted to a single scalar denoted by Z and a single fermion labeled by ψ . Just as in the $\mathfrak{su}(2)$ sector, we use the fermionic position-space oscillators b_j^\dagger, b_j to create

or annihilate fermionic ψ insertions in a ground state composed of K scalars:

$$|K\rangle = \text{tr}(Z^K) , \quad b_j^\dagger |K\rangle = \text{tr}(Z_1 \cdots Z_{j-1} \psi Z_{j+1} \cdots Z_K) . \quad (2.2.1)$$

In [33], Beisert gave the action of the Hamiltonian on the $\mathfrak{su}(2|3)$ spin chain to three-loop order.⁴ In the notation of [33], the action of the Hamiltonian on basis states can be represented in terms of special permutation operators denoted by

$$\left\{ \begin{matrix} A_1 \dots A_N \\ B_1 \dots B_N \end{matrix} \right\} ,$$

which replace all occurrences of the upper sequence of fields $A_1 \dots A_N$ in the trace by the lower sequence $B_1 \dots B_N$. Restricting Beisert's $\mathfrak{su}(2|3)$ Hamiltonian to the $\mathfrak{su}(1|1)$ subsector at one-loop order yields

$$H_{\mathfrak{su}(1|1)}^{(2)} = \left\{ \begin{matrix} Z\psi \\ Z\psi \end{matrix} \right\} + \left\{ \begin{matrix} \psi Z \\ \psi Z \end{matrix} \right\} - \left\{ \begin{matrix} Z\psi \\ \psi Z \end{matrix} \right\} - \left\{ \begin{matrix} \psi Z \\ Z\psi \end{matrix} \right\} + 2 \left\{ \begin{matrix} \psi\psi \\ \psi\psi \end{matrix} \right\} . \quad (2.2.2)$$

In terms of the position-space oscillators of eqn. (2.2.1), the $\mathfrak{su}(1|1)$ Hamiltonian can be assembled by inspection and takes the form

$$H_{\mathfrak{su}(1|1)}^{(2)} = \sum_{j=1}^K \left(b_j^\dagger b_j + b_{j+1}^\dagger b_{j+1} - b_{j+1}^\dagger b_j - b_j^\dagger b_{j+1} \right) . \quad (2.2.3)$$

There are no higher-body interaction terms at this order in λ . This fact can be checked by computing

$$\langle K | b_{i+1} b_i (H_{\mathfrak{su}(1|1)}^{(2)}) b_i^\dagger b_{i+1}^\dagger | K \rangle = 2 , \quad (2.2.4)$$

which reproduces the two-body matrix element given by the last term in eqn. (2.2.2).

⁴Beisert's three-loop Hamiltonian was restricted in [33] to the bosonic sector, but the author has since provided us with the complete version.

In momentum space we obtain

$$H_{\text{su}(1|1)}^{(2)} = 4 \sum_{p=0}^{K-1} \sin^2 \left(\frac{p\pi}{K} \right) \tilde{b}_p^\dagger \tilde{b}_p . \quad (2.2.5)$$

The two-loop $\text{su}(1|1)$ momentum-space Hamiltonian can be extracted in the same manner (the position-space version is too long to print here):

$$\begin{aligned} H_{\text{su}(1|1)}^{(4)} = & -8 \sum_{p=0}^{K-1} \sin^4 \left(\frac{p\pi}{K} \right) \tilde{b}_p^\dagger \tilde{b}_p + \frac{1}{4K} \sum_{p,q,r,s=0}^{K-1} \left\{ e^{\frac{2\pi i(q-2r)}{K}} + e^{\frac{2\pi i(2q-r)}{K}} - 4 e^{\frac{2\pi i(q-r)}{K}} \right. \\ & \left. - 2 e^{\frac{2\pi i(q-2r-s)}{K}} - 2 e^{\frac{2\pi i(q+s)}{K}} + e^{\frac{2\pi i(q-r+s)}{K}} + e^{\frac{2\pi i(2q-2r-s)}{K}} \right\} \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r \tilde{b}_s \delta_{p+q,r+s} . \end{aligned} \quad (2.2.6)$$

Finally, the complete three-loop Hamiltonian for this subsector is

$$\begin{aligned} H_{\text{su}(1|1)}^{(6)} = & 32 \sum_{p=0}^{K-1} \sin^6 \left(\frac{p\pi}{K} \right) \tilde{b}_p^\dagger \tilde{b}_p - \frac{1}{16} \sum_{p,q,r,s=0}^{K-1} e^{\frac{60\pi i(q-r)}{K}} \left\{ 2 e^{-\frac{2\pi i(27q-29r)}{K}} \right. \\ & + 2 e^{-\frac{2\pi i(28q-29r)}{K}} - 4 e^{-\frac{2\pi i(27q-28r)}{K}} + 37 e^{-\frac{2\pi i(29q-28r)}{K}} - 6 e^{-\frac{2\pi i(29q-27r)}{K}} \\ & + 8 e^{-\frac{56\pi i(q-r)}{K}} - 72 e^{-\frac{58\pi i(q-r)}{K}} - 6 e^{-\frac{2\pi i(29q-29r-2s)}{K}} - 40 e^{-\frac{2\pi i(29q-30r-s)}{K}} \\ & + 37 e^{-\frac{2\pi i(29q-29r-s)}{K}} - 8 e^{-\frac{2\pi i(29q-28r-s)}{K}} + 8 e^{-\frac{2\pi i(27q-28r+s)}{K}} + 2 e^{-\frac{2\pi i(28q-28r+s)}{K}} \\ & - 40 e^{-\frac{2\pi i(29q-28r+s)}{K}} - 4 e^{-\frac{2\pi i(27q-27r+s)}{K}} + 8 e^{-\frac{2\pi i(29q-27r+s)}{K}} + 2 e^{-\frac{2\pi i(27q-27r+2s)}{K}} \\ & \left. + 8 e^{-\frac{2\pi i(29q-30r-2s)}{K}} \right\} \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r \tilde{b}_s \delta_{p+q,r+s} + \frac{1}{16} \sum_{p,q,r,s,t,u=0}^{K-1} \left\{ 2 e^{\frac{2\pi i(q+2r-3s-2t)}{K}} \right. \\ & - e^{\frac{2\pi i(q+3r-3s-2t)}{K}} - 4 e^{\frac{2\pi i(q+2r-3s-t)}{K}} - e^{\frac{2\pi i(2q+3r-3s-t)}{K}} + 8 e^{\frac{2\pi i(q+2r-2s-t)}{K}} \\ & + 2 e^{\frac{2\pi i(2q+3r-2s-t)}{K}} - 4 e^{\frac{2\pi i(q+2r-3s-2t-u)}{K}} + 2 e^{\frac{2\pi i(q+3r-3s-2t-u)}{K}} + 2 e^{\frac{2\pi i(q+2r-2s+u)}{K}} \\ & \left. - 4 e^{\frac{2\pi i(q+2r-s+u)}{K}} - 4 e^{\frac{2\pi i(q+2r-2s-t+u)}{K}} \right\} \tilde{b}_p^\dagger \tilde{b}_q^\dagger \tilde{b}_r^\dagger \tilde{b}_s \tilde{b}_t \tilde{b}_u \delta_{p+q+r,s+t+u} . \end{aligned} \quad (2.2.7)$$

We note that $H_{\text{su}(1|1)}^{(2)}$, $H_{\text{su}(1|1)}^{(4)}$ and $H_{\text{su}(1|1)}^{(6)}$ terminate at one-body, two-body and three-body interactions, respectively. This will permit us to obtain the exact K -dependence of successive terms in the λ expansion of energy eigenvalues.

As in the $\mathfrak{su}(2)$ sector, we can use non-degenerate perturbation theory to extract the K^{-1} scaling coefficients of the $\mathfrak{su}(1|1)$ eigenvalue spectrum up to three loops in λ . The scaling coefficients extrapolated from numerical diagonalization of lattices up to $K = 40$ are recorded for one-loop, two-loop and three-loop orders in tables 2.5, 2.6 and 2.7, respectively. The same increase in leading power of K^{-1} with corresponding order in λ that was noted in the $\mathfrak{su}(2)$ sector is found here as well (we use the same notation for the scaling coefficients as before in order to keep track of these powers). It should also be noted that, because the impurities in this sector are fermions symmetrized on all group indices, the lattice momenta of all pseudoparticles must be different. These results amount to the following predictions for the one-loop and two-loop scaling coefficients:

$$\begin{aligned} E_{\mathfrak{su}(1|1)}^{(1,2)} &= (k_1^2 + k_1 k_2 + k_2^2) , & E_{\mathfrak{su}(1|1)}^{(1,3)} &= 0 , \\ E_{\mathfrak{su}(1|1)}^{(2,4)} - \frac{1}{4}(k_1^2 + k_1 k_2 + k_2^2)^2 , & & E_{\mathfrak{su}(1|1)}^{(2,5)} &= -(k_1^2 + k_1 k_2 + k_2^2)^2 . \end{aligned} \quad (2.2.8)$$

We again have the usual caveat that data on larger and larger lattices are required to maintain a fixed precision as one goes to higher and higher energy levels.

$E_{\mathfrak{su}(1 1)}^{(1,2)}$	$E_{\mathfrak{su}(1 1)}^{(1,3)}$	$E_{\mathfrak{su}(1 1)}^{(1,3)}/E_{\mathfrak{su}(1 1)}^{(1,2)}$	(k_1, k_2, k_3)
$1 + 1.3 \times 10^{-10}$	-1.9×10^{-8}	-1.9×10^{-8}	$(1, 0, -1)$
$4 - 1.0 \times 10^{-7}$	1.8×10^{-5}	4.6×10^{-6}	$(2, 0, -2)$
$7 - 2.5 \times 10^{-7}$	4.4×10^{-5}	6.3×10^{-6}	$(1, 2, -3)$
$7 - 2.5 \times 10^{-7}$	4.4×10^{-5}	6.3×10^{-6}	$(-1, -2, 3)$
$9 - 3.9 \times 10^{-7}$	7.9×10^{-5}	8.7×10^{-6}	$(3, 0, -3)$
$13 - 4.0 \times 10^{-6}$	8.2×10^{-4}	6.3×10^{-5}	$(1, 3, -4)$
$13 - 4.0 \times 10^{-6}$	8.2×10^{-4}	6.3×10^{-5}	$(-1, -3, 4)$
$16 - 2.0 \times 10^{-5}$	4.1×10^{-3}	2.6×10^{-4}	$(4, 0, -4)$
$19 - 3.5 \times 10^{-5}$	7.3×10^{-3}	3.8×10^{-4}	$(2, 3, -5)$
$19 - 3.5 \times 10^{-5}$	7.3×10^{-3}	3.8×10^{-4}	$(-2, -3, 5)$

Table 2.5: Scaling limit of one-loop numerical spectrum of three-impurity $\mathfrak{su}(1|1)$ subsector

$E_{\mathfrak{su}(1 1)}^{(2,4)}$	$E_{\mathfrak{su}(1 1)}^{(2,5)}$	$E_{\mathfrak{su}(1 1)}^{(2,5)}/E_{\mathfrak{su}(1 1)}^{(2,4)}$	(k_1, k_2, k_3)
-0.25	-0.99999	3.99995	(1, 0, -1)
-4.00006	-15.990	3.998	(2, 0, -2)
-12.251	-48.899	3.992	(1, 2, -3)
-12.251	-48.899	3.992	(-1, -2, 3)
-20.25	-80.89	3.995	(3, 0, -3)
-42.25	-168.2	3.98	(1, 3, -4)
-42.25	-168.2	3.98	(-1, -3, 4)
-64.00	-254.6	3.98	(4, 0, -4)
-90.26	-359.3	3.98	(2, 3, -5)
-90.26	-359.8	3.99	(-2, -3, 5)

Table 2.6: Scaling limit of two-loop numerical spectrum of three-impurity $\mathfrak{su}(1|1)$ subsector

The scaling limit of the three-loop ratio $E_{\mathfrak{su}(1|1)}^{(3,7)}/E_{\mathfrak{su}(1|1)}^{(3,6)}$ is recorded for the first few low-lying states in the spectrum in table 2.7.

The extrapolated gauge theory results in eqn. (2.2.8) for the one-loop coefficients $E_{\mathfrak{su}(1|1)}^{(1,3)}$ and $E_{\mathfrak{su}(1|1)}^{(1,2)}$ should be checked against the predictions of the general one-loop Bethe ansatz [32, 39] applied to the $\mathfrak{su}(1|1)$ sector (as far as we know, no higher-loop Bethe ansatz is available here). To apply the general Bethe ansatz equation of eqn. (2.1.29), we note that the $\mathfrak{su}(1|1)$ Dynkin diagram is just a single fermionic node: the Cartan matrix is empty and the single Dynkin label is $V_{\mathfrak{su}(1|1)} = 1$ [74, 75]. We therefore obtain the simple one-loop Bethe equation

$$\left(\frac{u_i + \frac{i}{2}}{u_i - \frac{i}{2}} \right)^K = 1 . \quad (2.2.9)$$

Rather remarkably, eqn. (2.2.9) can be solved exactly for arbitrary impurity number! The general $\mathfrak{su}(1|1)$ Bethe roots are

$$u_i = \frac{1}{2} \cot \left(\frac{k_i \pi}{K} \right) , \quad (2.2.10)$$

$E_{\mathfrak{su}(1 1)}^{(3,7)}/E_{\mathfrak{su}(1 1)}^{(3,6)}$	(k_1, k_2, k_3)
-86.41	(1, 0, -1)
-85.71	(2, 0, -2)
-83.74	(1, 2, -3)
-83.74	(-1, -2, 3)
-101.9	(3, 0, -3)
-96.01	(1, 3, -4)
-96.01	(-1, -3, 4)
-158.1	(4, 0, -4)

Table 2.7: Scaling limit of three-loop numerical spectrum of three-impurity $\mathfrak{su}(1|1)$ fermionic subsector

and the energy eigenvalues computed from eqn. (2.1.31) are

$$E_{\mathfrak{su}(1|1)} = 4 \sum_{i=1}^I \sin^2 \left(\frac{\pi k_i}{K} \right), \quad (2.2.11)$$

with the usual condition $\sum k_i = 0 \pmod K$ from eqn. (2.1.30). This is just the sum of free lattice Laplacian energies and clearly matches the energies one would obtain from the one-loop $\mathfrak{su}(1|1)$ Hamiltonian of eqn. (2.2.5) (since the latter has no interaction terms). No expansion in $1/K$ was necessary in this argument, but it is straightforward to expand the energies in $1/K$ and verify the numerical results obtained in table 2.5 and eqn. (2.2.8).

2.3 The $\mathfrak{sl}(2)$ sector

As noted in [38], integrable $\mathfrak{sl}(2)$ spin chains have previously been the subject of several studies involving, among other interesting problems, high-energy scattering amplitudes in non-supersymmetric QCD (see, e.g., [77] and references therein). The $\mathfrak{sl}(2)$ closed sector of $\mathcal{N} = 4$ SYM was studied in [69], and the spin chain Hamiltonian in this sector is presently known to one loop in λ . (For more recent progress, see ref. [52].)

The constituent fields in this sector are $SO(6)$ bosons, Z , carrying a single unit of R -charge ($Z = \phi_5 + i\phi_6$ or, in the language of Chapter 1, $\phi^{\frac{11}{2}}$), and each lattice site on the $\mathfrak{sl}(2)$ spin chain is occupied by a single Z field acted on by any number of the spacetime covariant derivatives $\nabla \equiv \nabla_1 + i\nabla_2$. The total R -charge of a particular operator is therefore equal to the lattice length K , and an I -impurity operator basis is spanned by single-trace operators carrying all possible distributions of I derivatives among the K lattice sites:

$$\begin{aligned} & \text{Tr} (\nabla^I Z Z^{K-1}) \ , \quad \text{Tr} (\nabla^{I-1} Z \nabla Z Z^{K-2}) \ , \\ & \text{Tr} (\nabla^{I-1} Z Z \nabla Z Z^{K-3}) \ , \quad \dots \end{aligned} \quad (2.3.1)$$

The integer I counts the total number of derivatives in the operator and, since any number of impurities can occupy the same lattice site, one can think of n derivative insertions at the i^{th} lattice site as n bosonic oscillator excitations at the i^{th} lattice position:

$$(a_i^\dagger)^n |K\rangle \sim \text{Tr} (Z^{i-1} \nabla^n Z Z^{K-i}) \ , \dots \quad (2.3.2)$$

The ground state $|K\rangle$ is represented by a length K chain with no derivative insertions: $|K\rangle = \text{Tr} (Z^K)$.

The one-loop $\mathfrak{sl}(2)$ spin chain Hamiltonian (corresponding to the dilatation operator in this sector) was constructed in [69] and was defined by its action on basis states rather than directly expressed as an operator:

$$\begin{aligned} H_{\mathfrak{sl}(2)}^{(2)} &= \sum_{j=1}^K H_{j,j+1}^{\mathfrak{sl}(2)} \ , \\ H_{1,2}^{\mathfrak{sl}(2)} (a_1^\dagger)^j (a_2^\dagger)^{n-j} |K\rangle &= \sum_{j'=0}^n \left[\delta_{j=j'} (h(j) + h(n-j)) - \frac{\delta_{j \neq j'}}{|j-j'|} \right] (a_1^\dagger)^{j'} (a_2^\dagger)^{n-j'} |K\rangle \end{aligned} \quad (2.3.3)$$

(where $h(n) = 1 + \dots + 1/n$ are the harmonic numbers). In other words, $H_{\mathfrak{sl}(2)}^{(2)}$ is a sum over the position-space Hamiltonian $H_{j,j+1}^{\mathfrak{sl}(2)}$, which acts on the j^{th} and $(j+1)^{\text{th}}$ (neighboring) lattice sites; the action of $H_{j,j+1}^{\mathfrak{sl}(2)}$ can be summarized by the explicit form given for $H_{1,2}^{\mathfrak{sl}(2)}$ above. Since it is only defined by its action on the state $(a_1^\dagger)^j (a_2^\dagger)^{n-j} |K\rangle$, it is difficult to immediately translate $H_{\mathfrak{sl}(2)}^{(2)}$ to momentum space. However, it is possible to expand it in powers of fields and use eqn. (2.3.3) to iteratively determine the expansion coefficients. The virial argument furthermore tells us that higher powers in the fields will determine higher powers of K^{-1} in the expansion of the energy. For our current purposes, it suffices to know the Hamiltonian expanded out to terms of fourth order in the fields and this truncation of the Hamiltonian can easily be constructed by inspection:

$$\begin{aligned} H_{\mathfrak{sl}(2)}^{(2)} = & - \sum_{j=1}^K \left[\left(a_{j+1}^\dagger - 2a_j^\dagger + a_{j-1}^\dagger \right) \left(a_j - \frac{1}{2} a_j^\dagger a_j^2 \right) \right. \\ & \left. + \frac{1}{4} \left(a_{j+1}^{\dagger 2} - 2a_j^{\dagger 2} + a_{j-1}^{\dagger 2} \right) a_j^2 \right] + \dots . \end{aligned} \quad (2.3.4)$$

Transformation to momentum space gives

$$\begin{aligned} H_{\mathfrak{sl}(2)}^{(2)} = & \sum_{p=0}^{K-1} 4 \sin^2 \frac{p\pi}{K} \tilde{a}_p^\dagger \tilde{a}_p \\ & + \frac{1}{K} \sum_{p,q,r,s=0}^{K-1} \delta_{p+q,r+s} \left(-\sin^2 \frac{p\pi}{K} - \sin^2 \frac{q\pi}{K} + \sin^2 \frac{(p+q)\pi}{K} \right) \tilde{a}_p^\dagger \tilde{a}_q^\dagger \tilde{a}_r \tilde{a}_s + \dots . \end{aligned} \quad (2.3.5)$$

This Hamiltonian acts on an I -impurity Fock space spanned by the generic states

$$\tilde{a}_{k_1}^\dagger \tilde{a}_{k_2}^\dagger \tilde{a}_{k_3}^\dagger \dots |K\rangle , \quad (2.3.6)$$

with lattice momenta labeled by $k_i = 0, \dots, K-1$, and subject to the constraint $\sum_i k_i = 0 \pmod K$. Numerically diagonalizing this Hamiltonian on a range of lattice sizes, we obtain data from which we extract the numerical predictions for the one-

loop coefficients $E_{\mathfrak{sl}(2)}^{(1,2)}$ and $E_{\mathfrak{sl}(2)}^{(1,3)}$ presented in table 2.8. We arrive at the following predictions for the scaling coefficients

$$\begin{aligned} E_{\mathfrak{sl}(2)}^{(1,2)} &= (k_1^2 + k_1 k_2 + k_2^2) , & E_{\mathfrak{sl}(2)}^{(1,3)} / E_{\mathfrak{sl}(2)}^{(1,2)} &= -2 , & k_1 &\neq k_2 \neq k_3 , \\ E_{\mathfrak{sl}(2)}^{(1,2)} &= 3n^2 , & E_{\mathfrak{sl}(2)}^{(1,3)} / E_{\mathfrak{sl}(2)}^{(1,2)} &= -7/3 , & k_1 &= k_2 = n, k_3 = -2n , \end{aligned} \quad (2.3.7)$$

and we can easily verify that the agreement with table 2.8 is excellent.

$E_{\mathfrak{sl}(2)}^{(1,2)}$	$E_{\mathfrak{sl}(2)}^{(1,3)}$	$E_{\mathfrak{sl}(2)}^{(1,3)} / E_{\mathfrak{sl}(2)}^{(1,2)}$	(k_1, k_2, k_3)
$1 + 1.2 \times 10^{-9}$	$-2 - 3.1 \times 10^{-7}$	$-2 - 3.1 \times 10^{-7}$	$(1, 0, -1)$
$3 - 7.6 \times 10^{-9}$	$-7 + 1.9 \times 10^{-6}$	$-7/3 + 6.3 \times 10^{-7}$	$(1, 1, -2)$
$3 - 7.6 \times 10^{-9}$	$-7 + 1.9 \times 10^{-6}$	$-7/3 + 6.3 \times 10^{-7}$	$(-1, -1, 2)$
$4 - 2.8 \times 10^{-7}$	$-8 + 6.9 \times 10^{-6}$	$-2 + 1.7 \times 10^{-6}$	$(2, 0, -2)$
$7 - 2.9 \times 10^{-7}$	$-14 + 7.1 \times 10^{-5}$	$-2 + 1.0 \times 10^{-5}$	$(1, 2, -3)$
$7 - 2.9 \times 10^{-7}$	$-14 + 7.1 \times 10^{-5}$	$-2 + 1.0 \times 10^{-5}$	$(-1, -2, 3)$
$9 - 4.1 \times 10^{-7}$	$-18 + 1.0 \times 10^{-4}$	$-2 + 1.0 \times 10^{-5}$	$(3, 0, -3)$
$12 + 8.4 \times 10^{-7}$	$-28 - 1.5 \times 10^{-4}$	$-7/3 - 1.2 \times 10^{-5}$	$(2, 2, -4)$
$12 + 8.4 \times 10^{-7}$	$-28 - 1.5 \times 10^{-4}$	$-7/3 - 1.2 \times 10^{-5}$	$(-2, -2, 4)$
$13 - 7.0 \times 10^{-6}$	$-26 + 1.7 \times 10^{-3}$	$-2 + 1.3 \times 10^{-4}$	$(1, 3, -4)$
$13 - 7.0 \times 10^{-6}$	$-26 + 1.7 \times 10^{-3}$	$-2 + 1.3 \times 10^{-4}$	$(-1, -3, 4)$
$16 - 1.4 \times 10^{-6}$	$-32 + 3.9 \times 10^{-4}$	$-2 + 2.4 \times 10^{-5}$	$(4, 0, -4)$
$19 - 7.5 \times 10^{-6}$	$-38 + 2.2 \times 10^{-3}$	$-2 + 1.1 \times 10^{-4}$	$(2, 3, -5)$
$19 - 7.5 \times 10^{-6}$	$-38 + 2.2 \times 10^{-3}$	$-2 + 1.1 \times 10^{-4}$	$(-2, -3, 5)$
$21 - 3.4 \times 10^{-6}$	$-42 + 8.8 \times 10^{-4}$	$-2 + 4.2 \times 10^{-5}$	$(1, 4, -5)$
$21 - 3.4 \times 10^{-6}$	$-42 + 8.8 \times 10^{-4}$	$-2 + 4.2 \times 10^{-5}$	$(-1, -4, 5)$

Table 2.8: Scaling limit of numerical spectrum of three-impurity $\mathfrak{sl}(2)$ sector at one loop

The extrapolated predictions can again be checked against those of the corresponding one-loop Bethe ansatz equations. In the $\mathfrak{sl}(2)$ sector the highest weight is $-1/2$: the Dynkin diagram therefore has coefficient $V_{\mathfrak{sl}(2)} = -1$ and the Cartan

matrix is $M_{\mathfrak{sl}(2)} = 2$. The Bethe equations (2.1.29, 2.1.30) thus reduce to

$$\left(\frac{u_i - i/2}{u_i + i/2} \right)^K = \prod_{j \neq i}^n \left(\frac{u_i - u_j + i}{u_i - u_j - i} \right), \quad (2.3.8)$$

$$1 = \prod_i^n \left(\frac{u_i - i/2}{u_i + i/2} \right). \quad (2.3.9)$$

Apart from a crucial minus sign, this is identical to the $\mathfrak{su}(2)$ Bethe equation (2.1.34).

In the absence of bound states, eqn. (2.3.8) is satisfied by the following Bethe roots:

$$\begin{aligned} u_1 &= -\frac{2(1+K)k_1^2 - (4+K)k_1k_2 - (4+K)k_2^2}{2\pi k_1(k_2^2 + k_1k_2 - 2k_1^2)} + O(K^{-1}), \\ u_2 &= -\frac{2(1+K)k_2^2 - (4+K)k_1k_2 - (4+K)k_1^2}{2\pi k_2(k_1^2 + k_1k_2 - 2k_2^2)} + O(K^{-1}), \\ u_3 &= -\frac{2(1+K)k_1^2 + (8+5K)k_1k_2 + 2(1+K)k_2^2}{2\pi(k_1+k_2)(2k_1+k_2)(k_1+2k_2)} + O(K^{-1}). \end{aligned} \quad (2.3.10)$$

Using eqn. (2.1.31), we obtain

$$\begin{aligned} E_{\mathfrak{sl}(2)}^{(2)}(k_1, k_2) &= \frac{\lambda}{K^3} (k_1^2 + k_1k_2 + k_2^2) (K - 2) + O(K^{-4}), \\ &\quad (k_1 \neq k_2 \neq k_3). \end{aligned} \quad (2.3.11)$$

For the bound state characterized by $k_1 = k_2 = n$ and $k_3 = -2n$, the Bethe roots are

$$\begin{aligned} u_1 &= \frac{7 - 3\sqrt{K} + 3K}{6\pi n} + O(K^{-1/2}), \\ u_2 &= \frac{7 + 3\sqrt{K} + 3K}{6\pi n} + O(K^{-1/2}), \\ u_3 &= -\frac{4 + 3K}{12\pi n} + O(K^{-1/2}), \end{aligned} \quad (2.3.12)$$

with spin chain energy

$$E_{\mathfrak{sl}(2)}^{(2)}(n) = \frac{\lambda n^2}{K^3} (3K - 7) + O(K^{-4}), \quad (k_1 = k_2 = n, k_3 = -2n). \quad (2.3.13)$$

These results again agree with the numerical results in table 2.8.

2.4 Discussion

In this chapter we have demonstrated that the virial expansion of the $\mathcal{N} = 4$ SYM spin chain Hamiltonian for small impurity number provides a simple and reliable method for computing exact anomalous dimensions of multi-impurity operators at small scalar R -charge (chain length) and estimating with great precision the near-BMN scaling behavior of these dimensions as the R -charge becomes large. The latter application, which is suited to direct comparison of gauge theory predictions with corresponding results on the string side of the AdS/CFT correspondence, works well for three-impurity operators to three-loop order in λ in the $\mathfrak{su}(2)$ sector (the order to which the $\mathfrak{su}(2)$ Hamiltonian is known definitively). Specifically, the numerical predictions from the virial approach for the near-BMN scaling coefficients ($E_{\mathfrak{su}(2)}^{(1,2)}$, $E_{\mathfrak{su}(2)}^{(1,3)}$, $E_{\mathfrak{su}(2)}^{(2,4)}$, $E_{\mathfrak{su}(2)}^{(2,5)}$, $E_{\mathfrak{su}(2)}^{(3,6)}$ and $E_{\mathfrak{su}(2)}^{(3,7)}$) match corresponding results from the $\mathfrak{su}(2)$ long-range Bethe ansatz to three-loop order, and will eventually be shown to agree with near-plane-wave string theory predictions to two loops (the disagreement with string theory at three loops is by now an expected outcome in these studies; this will also be demonstrated below). We also find convincing agreement near the BMN limit between the virial approach and the Bethe ansatz results at one-loop order in the closed $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ subsectors. As a side result we have found in the $\mathfrak{su}(1|1)$ sector an *exact* (in chain length) agreement between the Bethe ansatz and the virial expansion for one-loop operator dimensions with arbitrary impurity number (this was only possible because the Bethe equations can be solved exactly in this subsector for any number of impurities). There are currently no higher-loop Bethe ansätze for the $\mathfrak{su}(1|1)$ system, however, so in this sense our numerical predictions go beyond the current state of Bethe ansatz technology (see [78] for further developments of higher-loop gauge theory physics in non- $\mathfrak{su}(2)$ sectors). Recent progress in developing reliable Bethe equations in the $\mathfrak{sl}(2)$ sector beyond one-loop order has been made by Staudacher in [52]. It would be very interesting to find a general long-range Bethe equation appropriate for $\mathcal{N} = 4$ SYM at higher loop-order in λ , both for compar-

ison with string predictions and with the virial approach studied here. (For recent developments in this direction, see reference [63].)

Chapter 3

A curvature expansion of $AdS_5 \times S^5$

In order to address specifically stringy aspects of the duality, it is typically necessary to consider simplifying limits of the canonical $AdS_5 \times S^5$ background. As described above, Metsaev [25] showed that, in a certain plane-wave geometry supported by a constant RR flux, lightcone gauge worldsheet string theory reduces to a free theory with the novel feature that the worldsheet bosons and fermions acquire a mass. This solution was later shown to be a Penrose limit of the familiar $AdS_5 \times S^5$ supergravity solution [22], and describes the geometry near a null geodesic boosted around the equator of the S^5 subspace. The energies of Metsaev's free string theory are thus understood to be those of a string in the full $AdS_5 \times S^5$ space, in the limit that the states are boosted to large angular momentum about an equatorial circle in the S^5 . Corrections to the string spectrum that arise if the string is given a large, but finite, boost can be computed. Comparison of the resulting interacting spectrum with corrections (in inverse powers of the R -charge) to the dimensions of the corresponding gauge theory operators largely (but not completely) confirms expectations from AdS/CFT duality (see [26, 30] for discussion). The purpose of this chapter is to describe in fairly complete detail the methods used to obtain the results presented in [26] (but only outlined in that paper). Some aspects of the purely bosonic side of this problem were studied by Parnachev and Ryzhov [79]. Although we find no disagreement with them, our approach differs from theirs in certain respects, most notably in taking full account of supersymmetry.

The approach is to take the GS superstring action on $AdS_5 \times S^5$, constructed using the formalism of Cartan forms and superconnections on the $SU(2, 2|4)/(SO(4, 1) \times SO(5))$ coset superspace [80], expand it in powers of the background curvature and finally eliminate unphysical degrees of freedom by lightcone gauge quantization. We treat the resulting interaction Hamiltonian in first-order degenerate perturbation theory to find the first corrections to the highly-degenerate pp-wave spectrum. The complexity of the problem is such that we are forced to resort to symbolic manipulation programs to construct and diagonalize the perturbation matrix. In this chapter we give a proof of principle by applying our methods to the subspace of two-impurity excitations of the string. We show that the spectrum organizes itself into correct extended supersymmetry multiplets whose energies match well (if not perfectly) with what is known about gauge theory anomalous dimensions.

In Section 3.1 we introduce the problem by considering the bosonic sector of the theory alone. We comment on some interesting aspects of the theory that arise when restricting to the point-particle (or zero-mode) subsector. In Section 3.2 we review the construction of the GS superstring action on $AdS_5 \times S^5$ as a nonlinear sigma model on the $SU(2, 2|4)/(SO(4, 1) \times SO(5))$ coset superspace. In Sections 3.3 and 3.4 we perform a large-radius expansion on the relevant objects in the theory, and carry out the lightcone gauge reduction, thereby extracting explicit curvature corrections to the pp-wave Hamiltonian. Section 3.5 presents results on the curvature-corrected energy spectrum, further expanded to linear order in the modified 't Hooft coupling $\lambda' = g_{YM}^2 N_c / J^2$; results from corresponding gauge theory calculations (at one loop in $\lambda = g_{YM}^2 N_c$) are summarized and compared with the string theory. In Section 3.6 we extend the string theory analysis to higher orders in λ' , and compare results with what is known about gauge theory operator dimensions at higher-loop order.

3.1 Strings beyond the Penrose limit

To introduce the computation of finite- J corrections to the pp-wave string spectrum, we begin by discussing the construction of the lightcone gauge worldsheet Hamiltonian for the bosonic string in the full $AdS_5 \times S^5$ background. The problem is much more complicated when fermions are introduced, and we will take up that aspect of the calculation in a later section. A study of the purely bosonic problem gives us the opportunity to explain various strategic points in a simpler context.

In convenient global coordinates, the $AdS_5 \times S^5$ metric can be written in the form

$$ds^2 = \widehat{R}^2 (-\cosh^2 \rho \, dt^2 + d\rho^2 + \sinh^2 \rho \, d\Omega_3^2 + \cos^2 \theta \, d\phi^2 + d\theta^2 + \sin^2 \theta \, d\tilde{\Omega}_3^2) , \quad (3.1.1)$$

where \widehat{R} denotes the radius of both the sphere and the AdS space, and $d\Omega_3^2$, $d\tilde{\Omega}_3^2$ denote separate three-spheres. The coordinate ϕ is periodic with period 2π and, strictly speaking, so is the time coordinate t . In order to accommodate string dynamics, it is necessary to pass to the covering space in which time is *not* taken to be periodic. This geometry, supplemented by an RR field with N_c units of flux on the sphere, is a consistent, maximally supersymmetric type IIB superstring background, provided that $\widehat{R}^4 = g_s N_c (\alpha')^2$ (where g_s is the string coupling).

In its initial stages, development of the AdS/CFT correspondence focused on the supergravity approximation to string theory in $AdS^5 \times S^5$. Recently, attention has turned to the problem of evaluating truly stringy physics in this background and studying its match to gauge theory physics. The obstacles to such a program, of course, are the general difficulty of quantizing strings in curved geometries, and the particular problem of defining the superstring in the presence of RR background fields. As noted above, the string quantization problem is partly solved by looking at the dynamics of a string that has been boosted to lightlike momentum along some direction, or, equivalently, by quantizing the string in the background obtained by taking the Penrose limit of the original geometry using the lightlike geodesic corresponding to the boosted trajectory. The simplest choice is to boost along the equator of the

S^5 or, equivalently, to take the Penrose limit with respect to the lightlike geodesic $\phi = t$, $\rho = \theta = 0$ and to quantize the system in the appropriate lightcone gauge.

To quantize about the lightlike geodesic at $\rho = \theta = 0$, it is helpful to make the reparameterizations

$$\cosh \rho = \frac{1 + z^2/4}{1 - z^2/4}, \quad \cos \theta = \frac{1 - y^2/4}{1 + y^2/4}, \quad (3.1.2)$$

and work with the metric

$$ds^2 = \widehat{R}^2 \left[- \left(\frac{1 + \frac{1}{4}z^2}{1 - \frac{1}{4}z^2} \right)^2 dt^2 + \left(\frac{1 - \frac{1}{4}y^2}{1 + \frac{1}{4}y^2} \right)^2 d\phi^2 + \frac{dz_k dz_k}{(1 - \frac{1}{4}z^2)^2} + \frac{dy_{k'} dy_{k'}}{(1 + \frac{1}{4}y^2)^2} \right]. \quad (3.1.3)$$

The $SO(8)$ vectors spanning the eight directions transverse to the geodesic are broken into two $SO(4)$ subgroups parameterized by $z^2 = z_k z^k$ with $k = 1, \dots, 4$, and $y^2 = y_{k'} y^{k'}$ with $k' = 5, \dots, 8$. This form of the metric is well-suited for the present calculation: the spin connection, which will be important for the superstring action, turns out to have a simple functional form and the AdS_5 and S^5 subspaces appear nearly symmetrically. This metric has the full $SO(4, 2) \times SO(6)$ symmetry associated with $AdS_5 \times S^5$, but only the translation symmetries in t and ϕ and the $SO(4) \times SO(4)$ symmetry of the transverse coordinates remain manifest. The translation symmetries mean that string states have a conserved energy ω , conjugate to t , and a conserved (integer) angular momentum J , conjugate to ϕ . Boosting along the equatorial geodesic is equivalent to studying states with large J , and the lightcone Hamiltonian gives eigenvalues for $\omega - J$ in that limit. On the gauge theory side, the S^5 geometry is replaced by an $SO(6)$ R -symmetry, and J corresponds to the eigenvalue of an $SO(2)$ R -symmetry generator. The AdS/CFT correspondence implies that string energies in the boosted limit should match operator dimensions in the limit of large R -charge (a limit in which perturbative evaluation of operator dimensions becomes legitimate).

On dimensional grounds, taking the $J \rightarrow \infty$ limit on the string states is equivalent to taking the $\widehat{R} \rightarrow \infty$ limit on the metric (in the right coordinates). The coordinate redefinitions

$$t \rightarrow x^+ - \frac{x^-}{2\widehat{R}^2}, \quad \phi \rightarrow x^+ + \frac{x^-}{2\widehat{R}^2}, \quad z_k \rightarrow \frac{z_k}{\widehat{R}}, \quad y_{k'} \rightarrow \frac{y_{k'}}{\widehat{R}} \quad (3.1.4)$$

make it possible to take a smooth $\widehat{R} \rightarrow \infty$ limit. Expressing the metric (3.1.3) in these new coordinates, we obtain the following expansion in powers of $1/\widehat{R}^2$:

$$\begin{aligned} ds^2 \approx & 2 dx^+ dx^- + dz^2 + dy^2 - (z^2 + y^2) (dx^+)^2 \\ & + [2(z^2 - y^2) dx^- dx^+ + z^2 dz^2 - y^2 dy^2 - (z^4 - y^4) (dx^+)^2] \frac{1}{2\widehat{R}^2} \\ & + O(1/\widehat{R}^4). \end{aligned} \quad (3.1.5)$$

The leading \widehat{R} -independent part is the well-known pp-wave metric. The coordinate x^+ is dimensionless, x^- has dimensions of length squared, and the transverse coordinates now have dimensions of length. Since it is quadratic in the eight transverse bosonic coordinates, the pp-wave limit leads to a quadratic (and hence soluble) Hamiltonian for the bosonic string. The $1/\widehat{R}^2$ corrections to the metric are what will eventually concern us: they will add quartic interactions to the lightcone Hamiltonian and lead to first-order shifts in the energy spectrum of the string.

After introducing lightcone coordinates x^\pm according to (3.1.4), the general $AdS_5 \times S^5$ metric can be cast in the form

$$ds^2 = 2G_{+-} dx^+ dx^- + G_{++} dx^+ dx^+ + G_{--} dx^- dx^- + G_{AB} dx^A dx^B, \quad (3.1.6)$$

where x^A ($A = 1, \dots, 8$) labels the eight transverse directions, the metric components are functions of the x^A only, and the components G_{+A} and G_{-A} are not present. This simplifies even further for the pp-wave metric, where $G_{--} = 0$ and $G_{+-} = 1$. We will use (3.1.6) as the starting point for constructing the lightcone gauge worldsheet Hamiltonian (as a function of the transverse x^A and their conjugate momenta p_A)

and for discussing its expansion about the free pp-wave Hamiltonian.

The general bosonic Lagrangian density has a simple expression in terms of the target space metric:

$$\mathcal{L} = \frac{1}{2} h^{ab} G_{\mu\nu} \partial_a x^\mu \partial_b x^\nu , \quad (3.1.7)$$

where h is built out of the worldsheet metric γ according to $h^{ab} = \sqrt{-\det \gamma} \gamma^{ab}$ and the indices a, b label the worldsheet coordinates σ, τ . Since $\det h = -1$, there are only two independent components of h . The canonical momenta (and their inversion in terms of velocities) are

$$p_\mu = h^{\tau a} G_{\mu\nu} \partial_a x^\nu , \quad \dot{x}^\mu = \frac{1}{h^{\tau\tau}} G^{\mu\nu} p_\nu - \frac{h^{\tau\sigma}}{h^{\tau\tau}} x'^\mu . \quad (3.1.8)$$

The Hamiltonian density $H = p_\mu \dot{x}^\mu - \mathcal{L}$ is

$$H = \frac{1}{2h^{\tau\tau}} (p_\mu G^{\mu\nu} p_\nu + x'^\mu G_{\mu\nu} x'^\nu) - \frac{h^{\tau\sigma}}{h^{\tau\tau}} (x'^\mu p_\mu) . \quad (3.1.9)$$

As is usual in theories with general coordinate invariance (on the worldsheet in this case), the Hamiltonian is a sum of constraints times Lagrange multipliers built out of metric coefficients ($1/h^{\tau\tau}$ and $h^{\tau\sigma}/h^{\tau\tau}$).

One can think of the dynamical system we wish to solve as being defined by $\mathcal{L} = p_\mu \dot{x}^\mu - H$ (a phase space Lagrangian) regarded as a function of the coordinates x^μ , the momenta p_μ and the components h^{ab} of the worldsheet metric. To compute the quantum path integral, the exponential of the action constructed from this Lagrangian is functionally integrated over each of these variables. For a spacetime geometry like (3.1.6), one finds that with a suitable gauge choice for the worldsheet coordinates (τ, σ) , the functional integrations over all but the transverse (physical) coordinates and momenta can be performed, leaving an effective path integral for these physical variables. This is the essence of the lightcone approach to quantization.

The first step is to eliminate integrations over x^+ and p_- by imposing the lightcone

gauge conditions $x^+ = \tau$ and $p_- = \text{const.}$ (At this level of analysis, which is essentially classical, we will not be concerned with ghost determinants arising from this gauge choice.) As noted above, integrations over the worldsheet metric cause the coefficients $1/h^{\tau\tau}$ and $h^{\tau\sigma}/h^{\tau\tau}$ to act as Lagrange multipliers, generating delta functions that impose two constraints:

$$\begin{aligned} x'^- p_- + x'^A p_A &= 0 , \\ G^{++} p_+^2 + 2G^{+-} p_+ p_- + G^{--} p_-^2 + p_A G^{AB} p_B + x'^A G_{AB} x'^B + G_{--} \frac{(x'^A p_A)^2}{p_-^2} &= 0 . \end{aligned} \quad (3.1.10)$$

When integrations over x^- and p_+ are performed, the delta functions imposing constraints serve to evaluate x^- and p_+ in terms of the dynamical transverse variables (and the constant p_-). The first constraint is linear in x^- and yields $x'^- = -x'^A p_A / p_-$. Integrating this over σ and using the periodicity of x^- yields the standard level-matching constraint, without any modifications. The second constraint is quadratic in p_+ and can be solved explicitly for $p_+ = -H_{\text{LC}}(x^A, p_A)$. The remaining transverse coordinates and momenta have dynamics that follow from the phase space Lagrangian

$$\mathcal{L}_{\text{ps}} = p_+ + p_- \dot{x}^- + p_A \dot{x}^A \sim p_A \dot{x}^A - H_{\text{LC}}(x^A, p_A) , \quad (3.1.11)$$

where we have eliminated the p_- term by integrating by parts in time and imposing that p_- is constant. The essential result is that $-p_+ = H_{\text{LC}}$ is the Hamiltonian that generates evolution of the physical variables x^A , p_A in worldsheet time τ . This is, of course, dynamically consistent with the lightcone gauge identification $x^+ = \tau$ (which requires worldsheet and target space time translation to be the same).

We can solve the quadratic constraint equation (3.1.10) for $p_+ = -H_{\text{LC}}$ explicitly,

obtaining the uninspiring result

$$H_{\text{LC}} = -\frac{p_- G_{+-}}{G_{--}} - \frac{p_- \sqrt{G}}{G_{--}} \sqrt{1 + \frac{G_{--}}{p_-^2} (p_A G^{AB} p_B + x'^A G_{AB} x'^B) + \frac{G_{--}^2}{p_-^4} (x'^A p_A)^2} , \quad (3.1.12)$$

where

$$G \equiv G_{+-}^2 - G_{++} G_{--} . \quad (3.1.13)$$

This is not very useful as it stands, but we can put it in more manageable form by expanding it in powers of $1/\hat{R}^2$. We can actually do slightly better by observing that the constraint equation (3.1.10) becomes a linear equation for p_+ if $G_{--} = 0$ (which is equivalent to $G^{++} = 0$). Solving the linear equation for p_+ gives

$$H_{\text{LC}} = \frac{p_- G_{++}}{2G_{+-}} + \frac{G_{+-}}{2p_-} (p_A G^{AB} p_B + x'^A G_{AB} x'^B) , \quad (3.1.14)$$

a respectable non-linear sigma model Hamiltonian. In the general $AdS_5 \times S^5$ metric (3.1.1) we cannot define a convenient set of coordinates such that G_{--} identically vanishes. Using (3.1.4), however, we can find coordinates where G_{--} has an expansion which begins at $O(1/\hat{R}^4)$, while the other metric coefficients have terms of all orders in $1/\hat{R}^2$. Therefore, if we expand in $1/\hat{R}^2$ and keep terms of at most $O(1/\hat{R}^2)$, we may set $G_{--} = 0$ and use (3.1.14) to construct the expansion of the lightcone Hamiltonian to that order. The leading $O(\hat{R}^0)$ terms in the metric reproduce (as they should) the bosonic pp-wave Hamiltonian

$$H_{\text{LC}}^{\text{pp}} = \frac{1}{2} \left[(\dot{p}^A)^2 + (x'^A)^2 + (x^A)^2 \right] , \quad (3.1.15)$$

(choosing $p_- = 1$ for the conserved worldsheet momentum density). The $O(1/\hat{R}^2)$ terms generate a perturbing Hamiltonian density that is quartic in fields and quadratic

in worldsheet time and space derivatives:

$$H_{LC}^{\hat{R}^{-2}} = \frac{1}{4\hat{R}^2}(y^2 p_z^2 - z^2 p_y^2) + \frac{1}{4\hat{R}^2}((2z^2 - y^2)(z')^2 - (2y^2 - z^2)(y')^2) . \quad (3.1.16)$$

This is the bosonic part of the perturbing Hamiltonian we wish to derive. If we express it in terms of the creation and annihilation operators of the leading quadratic Hamiltonian (3.1.15) we can see that its matrix elements will be of order $1/J$, as will be the first-order perturbation theory shifts of the string energy eigenvalues. We defer the detailed discussion of this perturbation theory until we have the fermionic part of the problem in hand. Note that this discussion implies that if we wanted to determine the perturbed energies to higher orders in $1/\hat{R}^2$, we would have the very unpleasant problem of dealing with the square root form of the Hamiltonian (3.1.12).

We have to this point been discussing a perturbative approach to finding the effect of the true geometry of the $AdS_5 \times S^5$ background on the string spectrum. Before proceeding with this program, however, it is instructive to study a different limit in which the kinematics are unrestricted (no large- J limit is taken) but only modes of the string that are independent of the worldsheet coordinate (the zero-modes of the string) are kept in the Hamiltonian. This is the problem of quantizing the superparticle of the underlying supergravity in the $AdS_5 \times S^5$ background, a problem that has been solved many times (for references, see [81]). A remarkable fact, which seems not to have been explicitly observed before, is that the spectrum of the zero-mode Hamiltonian is *exactly* a sum of harmonic oscillators: the curvature corrections we propose to compute actually vanish on this special subspace. This fact is important to an understanding of the full problem, so we will make a brief digression to explain the solution to this toy problem.

The quantization of the superparticle in a supergravity background is equivalent to finding the eigensolutions of certain Laplacians, one for each spin that occurs in the superparticle massless multiplet. The point of interest to us can be made by analyzing the dynamics of the scalar particle and its associated scalar Laplacian, which only

depends on the background metric. With apologies, we will adopt another version of the $AdS_5 \times S^5$ metric, chosen because the scalar Laplacian is very simple in these coordinates:

$$ds^2 = -dt^2(\widehat{R}^2 + z^2) + d\phi^2(\widehat{R}^2 - y^2) + dz^j \left(\delta_{jk} - \frac{z^j z^k}{\widehat{R}^2 + z^2} \right) dz^k + dy^{j'} \left(\delta_{j'k'} + \frac{y^{j'} y^{k'}}{\widehat{R}^2 - y^2} \right) dy^{k'} . \quad (3.1.17)$$

As before, the coordinates z^k and $y^{k'}$ parameterize the two $SO(4)$ subspaces, and the indices j, k and j', k' run over $j, k = 1, \dots, 4$, and $j', k' = 5, \dots, 8$. This is a natural metric for analyzing fluctuations of a particle (or string) around the lightlike trajectory $\phi = t$ and $\vec{z} = \vec{y} = 0$. Because the metric components depend neither on t nor on ϕ , and because the problem is clearly separable in \vec{z} and \vec{y} , it makes sense to look for solutions of the form $\Phi = e^{-i\omega t} e^{iJ\phi} F(\vec{z}) G(\vec{y})$. The scalar Laplacian for ϕ in the above metric then reduces to

$$\left[-\frac{\omega^2}{\widehat{R}^2 + \vec{z}^2} + \frac{J^2}{\widehat{R}^2 - \vec{y}^2} - \frac{\partial}{\partial x^j} \left(\delta^{jk} + \frac{z^j z^k}{\widehat{R}^2} \right) \frac{\partial}{\partial z^k} - \frac{\partial}{\partial y^{j'}} \left(\delta^{j'k'} - \frac{y^{j'} y^{k'}}{\widehat{R}^2} \right) \frac{\partial}{\partial y^{k'}} \right] F(z) G(y) = 0 . \quad (3.1.18)$$

The radius \widehat{R} disappears from the equation upon rescaling the transverse coordinates by $z \rightarrow z/\widehat{R}$ and $y \rightarrow y/\widehat{R}$, so we can set $\widehat{R} = 1$ in what follows and use dimensional analysis to restore \widehat{R} if it is needed. The scalar Laplacian is essentially the lightcone Hamiltonian constraint (3.1.10) for string coordinates $z^k, y^{k'}$ and string momenta $p_z^k = -i\frac{\partial}{\partial z^k}$ and $p_y^{k'} = -i\frac{\partial}{\partial y^{k'}}$ (projected onto their zero modes). This implies that we can use the structure of the Laplacian to correctly order operators in the string Hamiltonian.

The periodicity $\phi \equiv \phi + 2\pi$ means that the angular momentum J is integrally quantized. The allowed values of ω then follow from the solution of the eigenvalue problem posed by (3.1.18). As the trial function Φ indicates, (3.1.18) breaks into

separate problems for \vec{z} and \vec{y} :

$$\begin{aligned} H_{AdS_5} F(\vec{z}) &= \left[p_j^z (\delta^{jk} + z^j z^k) p_k^z + \omega^2 \frac{z_k z^k}{1 + (z_k z^k)^2} \right] F(\vec{z}) = A(\omega) F(\vec{z}) , \\ H_{S^5} G(\vec{y}) &= \left[p_{j'}^y (\delta^{j'k'} - y^{j'} y^{k'}) p_{k'}^y + J^2 \frac{y_{k'} y^{k'}}{1 - (y_{k'} y^{k'})^2} \right] G(\vec{y}) = B(J) G(\vec{y}) , \end{aligned} \quad (3.1.19)$$

where $\omega^2 - J^2 = A + B$. The separation eigenvalues A, B depend on their respective parameters ω, J , and we determine the energy eigenvalues ω by finding the roots of the potentially complicated equation $\omega^2 - J^2 - A - B = 0$. The scalar Laplacian (3.1.18) is equivalent to the constraint equation (3.1.10) projected onto string zero modes, and we are once again seeing that the constraint doesn't directly give the Hamiltonian but rather an equation (quadratic or worse) to be solved for the Hamiltonian.

The H_{S^5} equation is just a repackaging of the problem of finding the eigenvalues of the $SO(6)$ Casimir invariant (another name for the scalar Laplacian on S^5) and H_{AdS_5} poses the corresponding problem for $SO(4, 2)$. The $SO(6)$ eigenvalues are obviously discrete, and the $SO(4, 2)$ problem also turns out to be discrete when one imposes the condition of finiteness at $z^2 \rightarrow \infty$ on the eigenfunctions (this is a natural restriction in the context of the AdS/CFT correspondence; for a detailed discussion see [81]). Thus we expect ω to have a purely discrete spectrum, with eigenvalues labeled by a set of integers. The simplest way to solve for the spectrum is to expand $F(\vec{z})$ and $G(\vec{y})$ in $SO(4)$ harmonics (since this symmetry is explicit), recognize that the radial equation is, in both cases, an example of Riemann's differential equation and then use known properties of the hypergeometric function to find the eigenvalues and eigenfunctions of (3.1.19). Since it takes three integers to specify an $SO(4)$ harmonic and one to specify a radial quantum number, we expect each of the two separated equations to have a spectrum labeled by four integers. The exact results for the separation eigenvalues turn out to be remarkably simple:

$$A = 2\omega \sum_1^4 \left(n_i + \frac{1}{2} \right) - \left[\sum_1^4 \left(n_i + \frac{1}{2} \right) \right]^2 + 4 , \quad n_i = 0, 1, 2, \dots ,$$

$$B = 2J \sum_1^4 \left(m_i + \frac{1}{2} \right) + \left[\sum_1^4 \left(m_i + \frac{1}{2} \right) \right]^2 + 4, \quad m_i = 0, 1, 2, \dots \quad (3.1.20)$$

Different eigenfunctions correspond to different choices of the collection of eight integers $\{n_i, m_i\}$, and the fact that the energies depend only on Σn_i and Σm_i correctly accounts for the degeneracy of eigenvalues. The special form of A and B means that the equation for the energy eigenvalue, $\omega^2 - J^2 - A - B = 0$, can be factored as

$$\begin{aligned} & \left[\omega - J - \sum_1^4 \left(n_i + \frac{1}{2} \right) - \sum_1^4 \left(m_i + \frac{1}{2} \right) \right] \\ & \times \left[\omega + J - \sum_1^4 \left(n_i + \frac{1}{2} \right) + \sum_1^4 \left(m_i + \frac{1}{2} \right) \right] = 0. \end{aligned} \quad (3.1.21)$$

For obvious reasons, we retain the root that assigns only positive values to ω , the energy conjugate to the global time t :

$$\omega - J = \sum_1^4 \left(n_i + \frac{1}{2} \right) + \sum_1^4 \left(m_i + \frac{1}{2} \right). \quad (3.1.22)$$

From the string point of view, ω catalogs the eigenvalues of the string worldsheet Hamiltonian restricted to the zero-mode subspace. Quite remarkably, it is an exact sum of harmonic oscillators, independent of whether J (and ω) are large or not. This is simply to say that the eigenvalues of the string Hamiltonian restricted to the zero-mode sector receive no curvature corrections and could have been calculated from the pp-wave string Hamiltonian (3.1.15). We have only shown this for the massless bosons of the theory, but we expect the same thing to be true for all the massless fields of type IIB supergravity. The implication for a perturbative account of the string spectrum is that states created using only zero-mode oscillators (of any type) will receive no curvature corrections. This feature will turn out to be a useful consistency check on

our quantization procedure. It is of course not true for a general classical background and is yet another manifestation of the special nature of the $AdS_5 \times S^5$ geometry.

3.2 GS superstring action on $AdS_5 \times S^5$

The $AdS_5 \times S^5$ target space can be realized as the coset superspace

$$G/H = \frac{SU(2, 2|4)}{SO(4, 1) \times SO(5)} . \quad (3.2.1)$$

The bosonic reduction of this coset is precisely $SO(4, 2) \times SO(6)/SO(4, 1) \times SO(5) \equiv AdS_5 \times S^5$. To quantize the theory, we will expand the action about a classical trajectory that happens to be invariant under the stabilizer group H . There is a general strategy for constructing a non-linear sigma model on a super-coset space in terms of the Cartan one-forms and superconnections of the super-coset manifold. In such a construction, the symmetries of the stabilizer subgroup remain manifest in the action while the remaining symmetries are nonlinearly realized (see, e.g., [24, 80, 82–85]). Metsaev and Tseytlin [83] carried out this construction for the $AdS_5 \times S^5$ geometry, producing a κ -symmetric, type IIB superstring action possessing the full $PSU(2, 2|4)$ supersymmetry of $AdS_5 \times S^5$. Their action is conceptually simple, comprising a kinetic term and a Wess–Zumino term built out of Cartan (super)one-forms on the super-coset manifold in the following way (this form was first presented in [86]):

$$\mathcal{S} = -\frac{1}{2} \int_{\partial M_3} d^2\sigma \, h^{ab} L_a^\mu L_b^\mu + i \int_{M_3} s^{IJ} L^\mu \wedge \bar{L}^I \Gamma^\mu \wedge L^J . \quad (3.2.2)$$

Repeated upper indices are summed over a Minkowskian inner product. The indices a, b are used to indicate the worldsheet coordinates (τ, σ) , and we use the values $a, b = 0$ to indicate the worldsheet time direction τ , and $a, b = 1$ to specify the σ direction. The matrix s^{IJ} is defined by $s^{IJ} \equiv \text{diag}(1, -1)$, where $I, J = 1, 2$. The Wess–Zumino term appears as an integral over a three-manifold M_3 , while the kinetic

term is integrated over the two-dimensional boundary ∂M_3 . The left-invariant Cartan forms are defined in terms of the coset space representative G by

$$G^{-1}dG = L^\mu P^\mu + L^\alpha \bar{Q}_\alpha + \bar{L}^\alpha Q_\alpha + \frac{1}{2}L^{\mu\nu} J^{\mu\nu} ,$$

$$L^N = dX^M L_M^N , \quad L_a^N = L_M^N \partial_a X^M , \quad X^M = (x^\mu, \theta^\alpha, \bar{\theta}^\alpha) . \quad (3.2.3)$$

The explicit expansion of this action in terms of independent fermionic degrees of freedom is rather intricate. One starts with two 32-component Majorana-Weyl spinors in ten dimensions: θ^I , where $I = 1, 2$ labels the two spinors. In a suitably chosen representation for the 32×32 ten-dimensional gamma matrices Γ^μ , the Weyl projection reduces to picking out the upper 16 components of θ and the surviving spinors can be combined into one complex 16-component spinor ψ :

$$\theta^I = \begin{pmatrix} \theta^\alpha \\ 0 \end{pmatrix}^I , \quad \psi^\alpha = \sqrt{2} [(\theta^\alpha)^1 + i(\theta^\alpha)^2] . \quad (3.2.4)$$

The following representation for Γ^μ (which has the desired property that $\Gamma_{11} = (\mathbf{1}_8, -\mathbf{1}_8)$) allows us to express their action on ψ in terms of real 16×16 γ -matrices:

$$\Gamma^\mu = \begin{pmatrix} 0 & \gamma^\mu \\ \bar{\gamma}^\mu & 0 \end{pmatrix} , \quad \gamma^\mu \bar{\gamma}^\nu + \gamma^\nu \bar{\gamma}^\mu = 2\eta^{\mu\nu} ,$$

$$\gamma^\mu = (1, \gamma^A, \gamma^9) , \quad \bar{\gamma}^\mu = (-1, \gamma^A, \gamma^9) . \quad (3.2.5)$$

The indices $\mu, \nu, \rho = 0, \dots, 9$ denote $SO(9, 1)$ vectors, and we will denote the corresponding spinor indices by $\alpha, \beta, \gamma, \delta = 1, \dots, 16$ (we also use the convention that upper-case indices $A, B, C, D = 1, \dots, 8$ indicate vectors of $SO(8)$, while $i, j, k = 1, \dots, 4$ ($i', j', k' = 5, \dots, 8$) indicate vectors from the $SO(3, 1) \cong SO(4)$ ($SO(4)$) subspaces associated with AdS_5 and S^5 respectively). The matrix γ^9 is formed by taking the product of the eight γ^A . A representation of γ^A matrices that will be convenient for explicit calculation is given in Appendix A. We also note that in the course of quantization we will impose the fermionic lightcone gauge fixing condition $\bar{\gamma}^9 \psi = \psi$.

This restricts the worldsheet fermions to lie in the 8_s representation of $SO(8)$ (and projects out the 8_c spinor), thus reducing the number of independent components of the worldsheet spinor from 16 to 8. The symmetric matrix

$$\Pi \equiv \gamma^1 \bar{\gamma}^2 \gamma^3 \bar{\gamma}^4 \quad (3.2.6)$$

appears in a number of places in the expansion of the action, so we give it an explicit definition. Since $\Pi^2 = 1$, it has eigenvalues ± 1 , which turn out to provide a useful sub-classification of the 8 components of the 8_s worldsheet spinor into two groups of 4. The quantity $\tilde{\Pi} = \Pi \gamma_9$ also appears, but does not require a separate definition because $\Pi \psi = \tilde{\Pi} \psi$ for spinors satisfying the lightcone gauge restriction to the 8_s representation.

Kallosh, Rahmfeld and Rajaraman presented in [82] a general solution to the supergravity constraints (Maurer-Cartan equations) for coset spaces exhibiting a superconformal isometry algebra of the form

$$\begin{aligned} [B_\mu, B_\nu] &= f_{\mu\nu}^\rho B_\rho, \\ [F_\alpha, B_\nu] &= f_{\alpha\nu}^\beta F_\beta, \\ \{F_\alpha, F_\beta\} &= f_{\alpha\beta}^\mu B_\mu, \end{aligned} \quad (3.2.7)$$

with B_μ and F_α representing bosonic and fermionic generators, respectively. In terms of these generators, the Cartan forms L^μ and superconnections L^α are determined completely by the structure constants $f_{\alpha\mu}^J$ and $f_{\alpha\beta}^\mu$:

$$L_{at}^\alpha = \left(\frac{\sinh t\mathcal{M}}{\mathcal{M}} \right)_\beta^\alpha (\mathcal{D}_a \theta)^\beta, \quad (3.2.8)$$

$$L_{at}^\mu = e^\mu{}_\nu \partial_a x^\nu + 2\theta^\alpha f_{\alpha\beta}^\mu \left(\frac{\sinh^2(t\mathcal{M}/2)}{\mathcal{M}^2} \right)_\gamma^\beta (\mathcal{D}_a \theta)^\gamma, \quad (3.2.9)$$

$$(\mathcal{M}^2)_\beta^\alpha = -\theta^\gamma f_{\gamma\mu}^\alpha \theta^\delta f_{\delta\beta}^\mu. \quad (3.2.10)$$

The dimensionless parameter t is used here to define “shifted” Cartan forms and

superconnections where, for example, $L_a^\mu = L_{at}^\mu|_{t=1}$. In the case of $AdS_5 \times S^5$, the Lagrangian takes the form

$$\mathcal{L}_{\text{Kin}} = -\frac{1}{2}h^{ab}L_a^\mu L_b^\mu, \quad (3.2.11)$$

$$\mathcal{L}_{\text{WZ}} = -2i\epsilon^{ab} \int_0^1 dt L_{at}^\mu s^{IJ} \bar{\theta}^I \Gamma^\mu L_{bt}^J. \quad (3.2.12)$$

In the context of eqns. (3.2.8, 3.2.9), it will be useful to choose a manifestation of the spacetime metric that yields a compact form of the spin connection. The form appearing in eqn. (3.1.3) is well suited to this requirement; the AdS_5 and S^5 subspaces are represented in (3.1.3) nearly symmetrically, and the spin connection is relatively simple:

$$\begin{aligned} \omega^{tz_k}_t &= \frac{z_k}{1 - \frac{1}{4}z^2}, & \omega^{z_j z_k}_{z_j} &= \frac{\frac{1}{2}z_k}{1 - \frac{1}{4}z^2}, \\ \omega^{\phi y_{k'}}_\phi &= -\frac{y_{k'}}{1 + \frac{1}{4}y^2}, & \omega^{y_{j'} y_{k'}}_{y_{j'}} &= -\frac{\frac{1}{2}y_{k'}}{1 + \frac{1}{4}y^2}. \end{aligned} \quad (3.2.13)$$

Upon moving to the lightcone coordinate system in (3.1.4), the x^+ direction remains null ($G_{--} = 0$) to $O(1/\hat{R}^4)$ in this expansion.

By introducing dimensionless contraction parameters Λ and Ω [87], one may express the $AdS_5 \times S^5$ isometry algebra keeping lightcone directions explicit:

$$\begin{aligned} [P^+, P^k] &= \Lambda^2 \Omega^2 J^{+k}, & [P^+, P^{k'}] &= -\Lambda^2 \Omega^2 J^{+k'}, \\ [P^+, J^{+k}] &= -\Lambda^2 P^k, & [P^+, J^{+k'}] &= \Lambda^2 P^{k'}, \\ [P^-, P^A] &= \Omega^2 J^{+A}, & [P^-, J^{+A}] &= P^A, \\ [P^j, P^k] &= \Lambda^2 \Omega^2 J^{jk}, & [P^{j'}, P^{k'}] &= -\Lambda^2 \Omega^2 J^{j'k'}, \\ [J^{+j}, J^{+k}] &= \Lambda^2 J^{jk}, & [J^{+j'}, J^{+k'}] &= -\Lambda^2 J^{j'k'}, \\ [P^j, J^{+k}] &= -\delta^{jk}(P^+ - \Lambda^2 P^-), & [P^r, J^{+s}] &= -\delta^{rs}(P^+ + \Lambda^2 P^-), \\ [P^i, J^{jk}] &= \delta^{ij}P^k - \delta^{ik}P^j, & [P^{i'}, J^{j'k'}] &= \delta^{i'j'}P^{k'} - \delta^{i'k'}P^{j'}, \\ [J^{+i}, J^{jk}] &= \delta^{ij}J^{+k} - \delta^{ik}J^{+j}, & [J^{+i'}, J^{j'k'}] &= \delta^{i'j'}J^{+k'} - \delta^{i'k'}J^{+j'}, \end{aligned}$$

$$[J^{ij}, J^{kl}] = \delta^{jk} J^{il} + 3 \text{ terms} , \quad [J^{i'j'}, J^{k'l'}] = \delta^{j'k'} J^{i'l'} + 3 \text{ terms} . \quad (3.2.14)$$

The bosonic sector of the algebra relevant to (3.2.7) takes the form

$$\begin{aligned} [J^{ij}, Q_\alpha] &= \frac{1}{2} Q_\beta (\gamma^{ij})^\beta_\alpha , \\ [J^{i'j'}, Q_\alpha] &= \frac{1}{2} Q_\beta (\gamma^{i'j'})^\beta_\alpha , \\ [J^{+i}, Q_\alpha] &= \frac{1}{2} Q_\beta (\gamma^{+i} - \Lambda^2 \gamma^{-i})^\beta_\alpha , \\ [J^{+i'}, Q_\alpha] &= \frac{1}{2} Q_\beta (\gamma^{+i'} + \Lambda^2 \gamma^{-i'})^\beta_\alpha , \\ [P^\mu, Q_\alpha] &= \frac{i\Omega}{2} Q_\beta (\Pi \gamma^+ \bar{\gamma}^\mu)^\beta_\alpha - \frac{i\Lambda^2 \Omega}{2} Q_\beta (\Pi \gamma^- \bar{\gamma}^\mu)^\beta_\alpha . \end{aligned} \quad (3.2.15)$$

The fermi-fermi anticommutation relations are

$$\begin{aligned} \{Q_\alpha, \bar{Q}_\beta\} &= -2i\gamma_{\alpha\beta}^\mu P^\mu - 2\Omega(\bar{\gamma}^k \Pi)_{\alpha\beta} J^{+k} - 2\Omega(\bar{\gamma}^{k'} \Pi)_{\alpha\beta} J^{+k'} \\ &\quad + \Omega(\bar{\gamma}^+ \gamma^{jk} \Pi)_{\alpha\beta} J^{jk} + \Omega(\bar{\gamma}^+ \gamma^{j'k'} \Pi)_{\alpha\beta} J^{j'k'} \\ &\quad - \Lambda^2 \Omega(\bar{\gamma}^- \gamma^{jk} \Pi)_{\alpha\beta} J^{jk} + \Lambda^2 \Omega(\bar{\gamma}^- \gamma^{j'k'} \Pi)_{\alpha\beta} J^{j'k'} . \end{aligned} \quad (3.2.16)$$

This form of the superalgebra has the virtue that one can easily identify the flat space ($\Omega \rightarrow 0$) and plane-wave ($\Lambda \rightarrow 0$) limits. The Maurer-Cartan equations in this coordinate system take the form

$$\begin{aligned} dL^\mu &= -L^{\mu\nu} L^\nu - 2i\bar{L}\bar{\gamma}^\mu L , \\ dL^\alpha &= -\frac{1}{4} L^{\mu\nu} (\gamma^{\mu\nu})^\alpha_\beta L^\beta + \frac{i\Omega}{2} L^\mu (\Pi \gamma^+ \bar{\gamma}^\mu)^\alpha_\beta L^\beta - \frac{i\Lambda^2 \Omega}{2} L^\mu (\Pi \gamma^- \bar{\gamma}^\mu)^\alpha_\beta L^\beta , \\ d\bar{L}^\alpha &= -\frac{1}{4} L^{\mu\nu} (\gamma^{\mu\nu})^\alpha_\beta \bar{L}^\beta - \frac{i\Omega}{2} L^\mu (\Pi \gamma^+ \bar{\gamma}^\mu)^\alpha_\beta \bar{L}^\beta + \frac{i\Lambda^2 \Omega}{2} L^\mu (\Pi \gamma^- \bar{\gamma}^\mu)^\alpha_\beta \bar{L}^\beta , \end{aligned} \quad (3.2.17)$$

where wedge products (3.2.2) are understood to be replaced by the following rules:

$$L^\mu L^\nu = -L^\nu L^\mu, \quad L^\mu L^\alpha = -L^\alpha L^\mu, \quad L^\alpha L^\beta = L^\beta L^\alpha. \quad (3.2.18)$$

Upon choosing a parameterization of the coset representative G

$$G(x, \theta) = f(x)g(\theta), \quad g(\theta) = \exp(\theta^\alpha \bar{Q}_\alpha + \bar{\theta}^\alpha Q_\alpha), \quad (3.2.19)$$

one derives a set of coupled differential equations for the shifted Cartan forms and superconnections:

$$\begin{aligned} \partial_t L_t &= d\theta + \frac{1}{4} L_t^{\mu\nu} \gamma^{\mu\nu} \theta - \frac{i\Omega}{2} L_t^\mu \Pi \gamma^+ \bar{\gamma}^\mu \theta + \frac{i\Lambda^2 \Omega}{2} L_t^\mu \Pi \gamma^- \bar{\gamma}^\mu \theta, \\ \partial_t L_t^\mu &= -2i\theta \bar{\gamma}^\mu \bar{L}_t - 2i\bar{\theta} \bar{\gamma}^\mu L_t, \\ \partial_t L_t^{-i} &= 2\Omega(\theta \bar{\gamma}^i \Pi \bar{L}_t) - 2\Omega(\bar{\theta} \bar{\gamma}^i \Pi L_t), \\ \partial_t L_t^{-r} &= 2\Omega(\theta \bar{\gamma}^r \Pi \bar{L}_t) - 2\Omega(\bar{\theta} \bar{\gamma}^r \Pi L_t), \\ \partial_t L_t^{ij} &= -2\Omega(\theta \bar{\gamma}^+ \gamma^{ij} \Pi \bar{L}_t) + 2\Omega(\bar{\theta} \bar{\gamma}^+ \gamma^{ij} \Pi L_t) \\ &\quad + 2\Lambda^2 \Omega(\theta \bar{\gamma}^- \gamma^{ij} \Pi \bar{L}_t) - 2\Lambda^2 \Omega(\bar{\theta} \bar{\gamma}^- \gamma^{ij} \Pi L_t), \\ \partial_t L_t^{i'j'} &= -2\Omega(\theta \bar{\gamma}^+ \gamma^{i'j'} \Pi \bar{L}_t) + 2\Omega(\bar{\theta} \bar{\gamma}^+ \gamma^{i'j'} \Pi L_t) - 2\Lambda^2 \Omega(\theta \bar{\gamma}^- \gamma^{i'j'} \Pi \bar{L}_t) \\ &\quad + 2\Lambda^2 \Omega(\bar{\theta} \bar{\gamma}^- \gamma^{i'j'} \Pi L_t). \end{aligned} \quad (3.2.20)$$

These coupled equations are subject to the following boundary conditions:

$$\begin{aligned} L_\pm(t=0) &= 0, \quad L_{t=0}^\mu = e^\mu, \quad L_{t=0}^\pm = e^\pm, \\ L_{t=0}^{\mu\nu} &= \omega^{\mu\nu}, \quad L_{t=0}^{-\mu} = \omega^{-\mu}. \end{aligned} \quad (3.2.21)$$

The generators $J^{-\mu}$ and $J^{kk'}$ are not present in the superalgebra, so the conditions

$$L^{+\mu} = 0, \quad L^{kk'} = 0 \quad (3.2.22)$$

are imposed as constraints.

To employ the general solution to the Maurer-Cartan equations (3.2.8, 3.2.9), the relevant sectors of the superalgebra may be rewritten in the more convenient 32-dimensional notation (setting $\Lambda = 1$ and $\Omega = 1$):

$$\begin{aligned} [Q_I, P^\mu] &= \frac{i}{2} \epsilon^{IJ} Q_J \Gamma_* \Gamma^\mu, \\ [Q_I, J^{\mu\nu}] &= -\frac{1}{2} Q_I \Gamma^{\mu\nu}, \\ \{(Q_I)^\mu, (Q_J)_\mu\} &= -2i\delta_{IJ} \Gamma^0 \Gamma^\rho P_\rho + \epsilon^{IJ} \left(-\Gamma^0 \Gamma^{jk} \Gamma_* J_{jk} + \Gamma^0 \Gamma^{j'k'} \Gamma'_* J_{j'k'} \right), \end{aligned} \quad (3.2.23)$$

where

$$\Gamma_* \equiv i\Gamma_{01234}, \quad \Gamma'_* \equiv i\Gamma_{56789}. \quad (3.2.24)$$

The Cartan forms and superconnections then take the following form:

$$\begin{aligned} L_{bt}^J &= \frac{\sinh t\mathcal{M}}{\mathcal{M}} \mathcal{D}_b \theta^J, \\ L_{at}^\mu &= e^\mu{}_\rho \partial_a x^\rho - 4i\bar{\theta}^I \Gamma^\mu \left(\frac{\sinh^2(t\mathcal{M}/2)}{\mathcal{M}^2} \right) \mathcal{D}_a \theta^I, \end{aligned} \quad (3.2.25)$$

where the covariant derivative is given by

$$(\mathcal{D}_a \theta)^I = \left(\partial_a \theta + \frac{1}{4} (\omega^{\mu\nu}{}_\rho \partial_a x^\rho) \Gamma^{\mu\nu} \theta \right)^I - \frac{i}{2} \epsilon^{IJ} e^\mu{}_\rho \partial_a x^\rho \Gamma_* \Gamma^\mu \theta^J. \quad (3.2.26)$$

The object \mathcal{M} is a 2×2 matrix, which, for convenience, is defined in terms of its square:

$$(\mathcal{M}^2)^{IL} = -\epsilon^{IJ} (\Gamma_* \Gamma^\mu \theta^J \bar{\theta}^L \Gamma^\mu) + \frac{1}{2} \epsilon^{KL} (-\Gamma^{jk} \theta^I \bar{\theta}^K \Gamma^{jk} \Gamma_* + \Gamma^{j'k'} \theta^I \bar{\theta}^K \Gamma^{j'k'} \Gamma'_*). \quad (3.2.27)$$

At this point, the GS action on $AdS_5 \times S^5$ (3.2.11, 3.2.12) may be expanded

to arbitrary order in fermionic and bosonic fields. In the present calculation, the parameters Ω and Λ remain set to unity, and the action is expanded in inverse powers of the target-space radius \widehat{R} , introduced in the rescaled lightcone coordinates in eqn. (3.1.4). The fact that supersymmetry must be protected at each order in the expansion determines a rescaling prescription for the fermions. Accordingly, the eight transverse bosonic directions x^A and the corresponding fermionic fields ψ^α receive a rescaling coefficient proportional to \widehat{R}^{-1} . The first curvature correction away from the plane-wave limit therefore occurs at quartic order in both bosonic and fermionic fluctuations. The particular lightcone coordinate system chosen in (3.1.4), however, gives rise to several complications. The x^\pm coordinates given by

$$t = x^+ - \frac{x^-}{2\widehat{R}^2} , \quad \phi = x^+ + \frac{x^-}{2\widehat{R}^2} \quad (3.2.28)$$

have conjugate momenta (in the language of BMN)

$$-p_+ = i\partial_{x^+} = i(\partial_t + \partial_\phi) = \Delta - J , \quad (3.2.29)$$

$$-p_- = i\partial_{x^-} = \frac{i}{2\widehat{R}^2}(\partial_\phi - \partial_t) = -\frac{1}{2\widehat{R}^2}(\Delta + J) , \quad (3.2.30)$$

with $\Delta = E = i\partial_t$ and $J = -i\partial_\phi$. The lightcone Hamiltonian is $H = -p_+$, so with $\Delta = J - p_+$ one may schematically write

$$\begin{aligned} p_- &= \frac{1}{2\widehat{R}^2}(2J - p_+) \\ &= \frac{J}{\widehat{R}^2} + \frac{H}{2\widehat{R}^2} \\ &= \frac{J}{\widehat{R}^2} \left(1 + \frac{1}{2J} \sum N\omega \right) . \end{aligned} \quad (3.2.31)$$

This result appears to be incorrect in the context of the lightcone gauge condition $\partial_\tau t = p_-$. To compensate for this, one must set the constant worldsheet density p_-

equal to something different from 1 (and non-constant) if the parameter length of the worldsheet is to be proportional to J . This operation introduces an additional $O(1/\widehat{R}^2)$ shift in the energy of the string oscillators. This is acceptable because, in practice, we wish to consider only degenerate subsets of energy states for comparison between the gauge theory and string theory results. Because of the compensation between corrections to J and the Hamiltonian contribution from p_- , the eigenvalues of J will remain constant within these degenerate subsets. Therefore, while it may seem incorrect to introduce operator-valued corrections to p_- , one could proceed pragmatically with the intent of restricting oneself to these degenerate subsets. When such a program is carried out, however, the resulting theory is subject to normal-ordering ambiguities; we instead use a coordinate system that is free of these complications.

A different choice of lightcone coordinates allows us to avoid this problem completely. By choosing

$$\begin{aligned} t &= x^+ , \\ \phi &= x^+ + \frac{x^-}{\widehat{R}^2} , \end{aligned} \tag{3.2.32}$$

we have

$$-p_+ = \Delta - J , \tag{3.2.33}$$

$$-p_- = i\partial_{x^-} = \frac{i}{\widehat{R}^2}\partial_\phi = -\frac{J}{\widehat{R}^2} , \tag{3.2.34}$$

such that p_- appears as a legitimate expansion parameter in the theory. In this coordinate system, the curvature expansion of the metric becomes

$$\begin{aligned} ds^2 &= 2dx^+dx^- - (x^A)^2(dx^+)^2 + (dx^A)^2 \\ &\quad + \frac{1}{\widehat{R}^2} \left[-2y^2dx^+dx^- + \frac{1}{2}(y^4 - z^4)(dx^+)^2 + (dx^-)^2 + \frac{1}{2}z^2dz^2 - \frac{1}{2}y^2dy^2 \right] \\ &\quad + O(\widehat{R}^{-4}) . \end{aligned} \tag{3.2.35}$$

The operator-valued terms in p_- that appear under the first coordinate choice (3.2.28) are no longer present. However, it will be shown that this new coordinate system induces correction terms to the spacetime curvature of the worldsheet metric. Furthermore, the appearance of a nonvanishing G_{--} component, and the loss of many convenient symmetries between terms associated with the x^+ and x^- directions bring some additional complications into the analysis. The advantage is that the results will be unambiguous in the end (and free from normal-ordering ambiguities).

3.3 Curvature corrections to the Penrose limit

In this section we expand the GS superstring action on $AdS_5 \times S^5$ in powers of $1/\widehat{R}^2$. We begin by constructing various quantities including combinations of Cartan one-forms relevant to the worldsheet Lagrangian. Spacetime curvature corrections to the worldsheet metric will be calculated by analyzing the x^- equation of motion and the covariant gauge constraints order-by-order.

We introduce the notation

$$\Delta_n^\mu \equiv \bar{\theta}^I \Gamma^\mu \mathcal{D}_0^n \theta^I, \quad (3.3.1)$$

$$\Delta'_n{}^\mu \equiv \bar{\theta}^I \Gamma^\mu \mathcal{D}_1^n \theta^I, \quad (3.3.2)$$

where the covariant derivative is expanded in powers of $(1/\widehat{R})$:

$$\mathcal{D}_a = \mathcal{D}_a^0 + \frac{1}{\widehat{R}} \mathcal{D}_a^1 + \frac{1}{\widehat{R}^2} \mathcal{D}_a^2 + O(\widehat{R}^{-3}). \quad (3.3.3)$$

Terms in the Wess-Zumino Lagrangian are encoded using a similar notation:

$$\square_n^\mu \equiv s^{IJ} \bar{\theta}^I \Gamma^\mu \mathcal{D}_0^n \theta^J, \quad (3.3.4)$$

$$\square'_n{}^\mu \equiv s^{IJ} \bar{\theta}^I \Gamma^\mu \mathcal{D}_1^n \theta^J. \quad (3.3.5)$$

The subscript notation $(\Delta_n^\mu)_{\theta^4}$ will be used to indicate the quartic fermionic term

involving \mathcal{M}^2 :

$$(\Delta_n^\mu)_{\theta^4} \equiv \frac{1}{12} \bar{\theta}^I (\mathcal{M}^2) \mathcal{D}_0^n \theta^I . \quad (3.3.6)$$

For the present, it will be convenient to remove an overall factor of \widehat{R}^2 from the definition of the vielbeins e^μ_ν . In practice, this choice makes it easier to recognize terms that contribute to the Hamiltonian at the order of interest, and, in the end, allows us to avoid imposing an additional rescaling operation on the fermions. We proceed by keeping terms to $O(1/\widehat{R}^4)$, with the understanding that an extra factor of \widehat{R}^2 must be removed in the final analysis. The covariant derivative

$$\mathcal{D}_a \theta^I = \partial_a \theta^I + \frac{1}{4} \partial_a x^\mu \omega_\mu^{\nu\rho} \Gamma_{\nu\rho} \theta^I - \frac{i}{2} \epsilon^{IJ} \Gamma_* \Gamma_\mu e^\mu_\nu \partial_a x^\nu \theta^J \quad (3.3.7)$$

may then be expanded to $O(1/\widehat{R}^2)$ (we will not need $O(1/\widehat{R}^3)$ terms, because the covariant derivative always appears left-multiplied by a spacetime spinor $\bar{\theta}$):

$$\begin{aligned} \mathcal{D}_0 \theta^I &= \left[\partial_0 \theta^I - p_- \epsilon^{IJ} \Pi \theta^J \right] + \frac{1}{\widehat{R}} \left[\frac{p_-}{4} \left(z_j \Gamma^{-j} - y_{j'} \Gamma^{-j'} \right) \theta^I + \frac{1}{4} \epsilon^{IJ} \Gamma^- \Pi (\dot{x}^A \Gamma^A) \theta^J \right] \\ &+ \frac{1}{\widehat{R}^2} \left[\frac{1}{4} (\dot{z}_j z_k \Gamma^{jk} - \dot{y}_{j'} y_{k'} \Gamma^{j'k'}) \theta^I + \frac{p_-}{4} \epsilon^{IJ} \Pi (y^2 - z^2) \theta^J \right. \\ &\left. - \frac{1}{2} \epsilon^{IJ} (\dot{x}^-) \Pi \theta^J \right] + O(\widehat{R}^{-3}) , \end{aligned} \quad (3.3.8)$$

$$\begin{aligned} \mathcal{D}_1 \theta^I &= \partial_1 \theta^I + \frac{1}{4\widehat{R}} \epsilon^{IJ} \Gamma^- \Pi (x'^A \Gamma^A) \theta^J \\ &+ \frac{1}{\widehat{R}^2} \left[\frac{1}{4} (z'_j z_k \Gamma^{jk} - y'_{j'} y_{k'} \Gamma^{j'k'}) \theta^I - \frac{1}{2} \epsilon^{IJ} (x'^-) \Pi \theta^J \right] + O(\widehat{R}^{-3}) . \end{aligned} \quad (3.3.9)$$

Note that we have not rescaled the spinor field θ in the above expansion. This allows us to isolate the bosonic scaling contribution from the covariant derivative when combining various terms in the Lagrangian. Subsequently, the fermionic rescaling is performed based on the number of spinors appearing in each term (two spinors for each Δ^μ or \square^μ , and four for each $(\Delta^\mu)_{\theta^4}$). The worldsheet derivative notation is given by $\partial_\tau x = \partial_0 x = \dot{x}$ and $\partial_\sigma x = \partial_1 x = x'$.

The various sectors of the worldsheet Lagrangian are assembled keeping x^- and its derivatives explicit; these will be removed by imposing the covariant gauge constraints. From the supervielbein and superconnection

$$\begin{aligned} L_{at}^\mu &= e^\mu{}_\nu \partial_a x^\nu - 4i\bar{\theta}^I \Gamma^\mu \left(\frac{\sinh^2(t\mathcal{M}/2)}{\mathcal{M}^2} \right) \mathcal{D}_a \theta^I \\ &\approx e^\mu{}_\nu \partial_a x^\nu - i\bar{\theta}^I \Gamma^\mu \left(t^2 + \frac{t^4 \mathcal{M}^2}{12} \right) \mathcal{D}_a \theta^I, \end{aligned} \quad (3.3.10)$$

$$L_{at}^I = \frac{\sinh t\mathcal{M}}{\mathcal{M}} \mathcal{D}_a \theta^I \approx \left(t + \frac{t^3}{6} \mathcal{M}^2 \right) \mathcal{D}_a \theta^I, \quad (3.3.11)$$

we form the following objects:

$$\begin{aligned} L_0^\mu L_0^\mu &= \frac{1}{\widehat{R}^2} \left\{ 2p_- \dot{x}^- - p_-^2 (x^A)^2 + (\dot{x}^A)^2 - 2ip_- \Delta_0^- \right\} \\ &+ \frac{1}{\widehat{R}^4} \left\{ (\dot{x}^-)^2 - 2p_- y^2 \dot{x}^- + \frac{1}{2} (\dot{z}^2 z^2 - \dot{y}^2 y^2) + \frac{p_-^2}{2} (y^4 - z^4) \right. \\ &- 2i \left[\frac{1}{2} \dot{x}^- \Delta_0^- + p_- \Delta_2^- + p_- (\Delta_0^-)_{\theta^4} \right. \\ &\left. \left. - \frac{p_-}{4} (y^2 - z^2) \Delta_0^- + \dot{x}^A \Delta_1^A \right] \right\} + O(\widehat{R}^{-6}), \end{aligned} \quad (3.3.12)$$

$$\begin{aligned} L_1^\mu L_1^\mu &= \frac{1}{\widehat{R}^2} (x'^A)^2 + \frac{1}{\widehat{R}^4} \left\{ \frac{1}{2} (z'^2 z^2 - y'^2 y^2) + (x'^-)^2 \right. \\ &\left. - 2ix'^A \Delta_1'^A - ix'^- \Delta_0'^- \right\} + O(\widehat{R}^{-6}), \end{aligned} \quad (3.3.13)$$

$$\begin{aligned} L_0^\mu L_1^\mu &= \frac{1}{\widehat{R}^2} \left\{ p_- x'^- + \dot{x}^A x'^A - ip_- \Delta_0'^- \right\} \\ &+ \frac{1}{\widehat{R}^4} \left\{ x'^- \dot{x}^- - p_- y^2 x'^- + \frac{1}{2} (z^2 \dot{z}_k z'_k - y^2 \dot{y}_{k'} y'_{k'}) \right. \\ &- ip_- \Delta_2'^- - ip_- (\Delta_0'^-)_{\theta^4} - i \frac{p_-}{4} (z^2 - y^2) \Delta_0'^- - \frac{i}{2} \dot{x}^- \Delta_0'^- - i \dot{x}^A \Delta_1'^A \\ &\left. - ix'^A \Delta_1^A - \frac{i}{2} x'^- \Delta_0^- \right\} + O(\widehat{R}^{-6}). \end{aligned} \quad (3.3.14)$$

It will be advantageous to enforce the lightcone gauge condition $x^+ = \tau$ at all orders in the theory.¹ When fermions are included, this choice allows us to keep the

¹ This differs from the approach presented in [79].

κ -symmetry condition $\Gamma^+\theta = 0$ exact. In the pp-wave limit, keeping the worldsheet metric flat in this lightcone gauge is consistent with the equations of motion. Beyond leading order, however, we are forced to consider curvature corrections to the worldsheet metric that appear in both the conformal gauge constraints and the worldsheet Hamiltonian. In the purely bosonic case described in Section 3.1 above, these corrections are kept implicit by defining gauge constraints in terms of canonical momenta. In the supersymmetric theory, we must explicitly calculate these corrections. The strategy is to expand the x^- equations of motion in rescaled coordinates (3.2.32) and solve for the components of the worldsheet metric order-by-order. By varying x^- in the full Lagrangian we obtain

$$\begin{aligned} \frac{\delta \mathcal{L}}{\delta \dot{x}^-} &= \frac{1}{2} h^{00} \left\{ \frac{2p_-}{\widehat{R}^2} + \frac{1}{\widehat{R}^4} [2\dot{x}^- - 2p_- y^2 - i\bar{\theta}^I \Gamma^- \partial_0 \theta^I + 2ip_- \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J] \right\} \\ &\quad + \frac{i}{2\widehat{R}^4} s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J + O(\widehat{R}^{-6}). \end{aligned} \quad (3.3.15)$$

The worldsheet metric is taken to be flat at leading order, so there is no contribution from $L_0^\mu L_1^\mu$ here. To obtain corrections to h^{ab} entirely in terms of physical variables, however, we must eliminate all instances of x^- (or its derivatives) from the above variation. We can solve the conformal gauge constraints at leading order to remove \dot{x}^- from (3.3.15). These constraints are obtained by varying the Lagrangian with respect to the worldsheet metric itself:

$$T_{ab} = L_a^\mu L_b^\mu - \frac{1}{2} h_{ab} h^{cd} L_c^\mu L_d^\mu, \quad (3.3.16)$$

yielding a symmetric traceless tensor with two independent components. To leading order in $1/\widehat{R}$, we find

$$\begin{aligned} T_{00} &= \frac{1}{2} (L_0^\mu L_0^\mu + L_1^\mu L_1^\mu) + \dots = 0 \\ &= \frac{1}{2\widehat{R}^2} \left(2p_- \dot{x}^- - p_-^2 (x^A)^2 + (\dot{x}^A)^2 - 2ip_- \Delta_0^- + (x'^A)^2 \right) + O(\widehat{R}^{-4}), \end{aligned}$$

$$\begin{aligned}
T_{01} &= L_0^\mu L_1^\mu + \cdots = 0 \\
&= p_- x'^- + \dot{x}^A x'^A - i p_- \Delta_0'^- + O(\widehat{R}^{-4}) .
\end{aligned} \tag{3.3.17}$$

Expanding \dot{x}^- and x'^- in the same fashion,

$$\dot{x}^- = \sum_n \frac{a_n}{\widehat{R}^n} , \quad x'^- = \sum_n \frac{a'_n}{\widehat{R}^n} , \tag{3.3.18}$$

we use (3.3.17) and (3.3.17) to obtain

$$a_0 = \frac{p_-}{2} (x^A)^2 - \frac{1}{2p_-} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] + i \bar{\theta}^I \Gamma^- \partial_0 \theta^I - i p_- \epsilon^{IJ} \bar{\theta}^I \Gamma^- \Pi \theta^J , \tag{3.3.19}$$

$$a'_0 = -\frac{1}{p_-} \dot{x}^A x'^A + i \bar{\theta}^I \Gamma^- \partial_1 \theta^I . \tag{3.3.20}$$

By substituting back into (3.3.15), and performing the analogous operation for the x'^- variation, these leading-order solutions provide the following expansions for the objects that enter into the x^- equation of motion:

$$\begin{aligned}
\frac{\delta \mathcal{L}}{\delta \dot{x}^-} &= \frac{1}{2} h^{00} \left\{ \frac{2p_-}{\widehat{R}^2} + \frac{1}{\widehat{R}^4} \left[p_- (z^2 - y^2) - \frac{1}{p_-} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] + i \bar{\theta}^I \Gamma^- \partial_0 \theta^I \right] \right\} \\
&\quad + \frac{i}{2 \widehat{R}^4} s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J + O(\widehat{R}^{-6}) , \\
\frac{\delta \mathcal{L}}{\delta x'^-} &= \frac{h^{01} p_-}{\widehat{R}^2} + \frac{h^{11}}{\widehat{R}^4} \left(-\frac{1}{p_-} \dot{x}^A x'^A + \frac{i}{2} \bar{\theta}^I \Gamma^- \partial_1 \theta^I \right) - \frac{i}{2 \widehat{R}^4} s^{IJ} \bar{\theta}^I \Gamma^- \partial_0 \theta^J + O(\widehat{R}^{-6}) .
\end{aligned} \tag{3.3.21}$$

It is obvious from these expressions that the x^- equation of motion will not be consistent with the standard choice of flat worldsheet metric ($h^{00} = -h^{11} = 1, h^{01} = 0$). We therefore expand h^{ab} in powers of \widehat{R}^{-1} , taking it to be flat at leading order and allowing the higher-order terms (the \tilde{h}^{ab}) to depend on the physical variables in some

way:

$$\begin{aligned} h^{00} &= -1 + \frac{\tilde{h}^{00}}{\widehat{R}^2} + O(\widehat{R}^{-4}) , & h^{11} &= 1 + \frac{\tilde{h}^{11}}{\widehat{R}^2} + O(\widehat{R}^{-4}) , \\ h^{01} &= \frac{\tilde{h}^{01}}{\widehat{R}^2} + O(\widehat{R}^{-4}) . \end{aligned} \quad (3.3.22)$$

Using (3.3.19) and (3.3.20), we find that the specific metric choice

$$\tilde{h}^{00} = \frac{1}{2}(z^2 - y^2) - \frac{1}{2p_-^2} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] + \frac{i}{2p_-} \bar{\theta}^I \Gamma^- \partial_0 \theta^I - \frac{i}{2p_-} s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J , \quad (3.3.23)$$

$$\tilde{h}^{01} = \frac{1}{p_-^2} \dot{x}^A x'^A - \frac{i}{2p_-} \bar{\theta}^I \Gamma^- \partial_1 \theta^I + \frac{i}{2p_-} s^{IJ} \bar{\theta}^I \Gamma^- \partial_0 \theta^J \quad (3.3.24)$$

simplifies the expressions of (3.3.21) to

$$\frac{\delta \mathcal{L}}{\delta \dot{x}^-} = 1 + O(\widehat{R}^{-4}) , \quad \frac{\delta \mathcal{L}}{\delta x'^-} = O(\widehat{R}^{-4}) . \quad (3.3.25)$$

The x^- equation of motion is then consistent with the standard lightcone gauge choice $\dot{x}^+ = p_-$ to $O(1/\widehat{R}^2)$ (with no corrections to p_- , which must remain constant). Note that $\tilde{h}^{00} = -\tilde{h}_{00}$ and $\tilde{h}_{00} = \tilde{h}_{11}$. The fact that these curvature corrections have bi-fermionic contributions is ultimately due to the presence of a non-vanishing G_{--} term in the expanded metric (3.2.35).

Since the worldsheet metric is known to $O(1/\widehat{R}^2)$, x^- can now be determined to this order from the covariant gauge constraints (3.3.16). By invoking the leading-order solutions (3.3.17, 3.3.17), we can simplify the equations to some extent:

$$T_{00} = \frac{1}{2} (L_0^\mu L_0^\mu + L_1^\mu L_1^\mu) + \frac{\tilde{h}^{00}}{\widehat{R}^2} L_1^\mu L_1^\mu + O(\widehat{R}^{-3}) = 0 , \quad (3.3.26)$$

$$T_{01} = L_0^\mu L_1^\mu - \frac{\tilde{h}_{01}}{\widehat{R}^2} L_1^\mu L_1^\mu + O(\widehat{R}^{-3}) = 0 . \quad (3.3.27)$$

Equation (3.3.26) may be expanded to solve for a_2 , the first subleading correction to

\dot{x}^- :

$$\begin{aligned}
T_{00} = & 2p_- a_2 + a_0^2 - 2p_- y^2 a_0 + a_0'^2 + \frac{1}{2}(\dot{z}^2 z^2 - \dot{y}^2 y^2) + \frac{p_-^2}{2}(y^4 - z^4) \\
& + \frac{1}{2}(z'^2 z^2 - y'^2 y^2) + (z^2 - y^2)(x'^A)^2 - \frac{1}{p_-^2} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] (x'^A)^2 \\
& + \frac{i}{p_-} (x'^A)^2 \bar{\theta}^I \Gamma^- \partial_0 \theta^I - \frac{i}{p_-} (x'^A)^2 s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J - i a_0 \Delta_0^- - 2i p_- \Delta_2^- \\
& - 2i p_- (\Delta_0^-)_{\theta^4} + \frac{i p_-}{2} (y^2 - z^2) \Delta_0^- - 2i (\dot{x}^A \Delta_1^A + x'^A \Delta_1^A) - i a_0' \Delta_0'^- = 0 .
\end{aligned} \tag{3.3.28}$$

The remaining independent component T_{01} is the current associated with translation symmetry on the closed-string worldsheet. Enforcing the constraint $T_{01} = 0$ is equivalent to imposing the level-matching condition on physical string states. This condition can be used to fix higher-order corrections to x'^- , as is required by conformal invariance on the worldsheet. However, since our goal is to examine curvature corrections to the pp-wave limit using first-order perturbation theory, we will only need to enforce the level-matching condition on string states that are eigenstates of the pp-wave theory. We therefore need only consider the equation $T_{01} = 0$ to leading order in the expansion, which yields (3.3.20) above. If we were interested in physical eigenstates of the geometry corrected to $O(1/\hat{R}^2)$ (i.e., solving the theory exactly to this order), we would be forced to solve $T_{01} = 0$ to $O(1/\hat{R}^2)$.

With solutions to the x^- equations of motion and an expansion of the worldsheet metric to the order of interest, we may proceed with expressing the Hamiltonian as the generator of lightcone time translation: $p_+ = \delta \mathcal{L} / \delta \dot{x}^+$. It is helpful to first vary Δ^μ with respect to $\partial_0 t$ and $\partial_0 \phi$:

$$\frac{\delta \Delta^\mu}{\delta(\partial_0 t)} = \bar{\theta}^I \Gamma^\mu \left[-\frac{1}{2\hat{R}^3} z_j \Gamma^{0j} \theta^I - \frac{1}{2} \epsilon^{IJ} \Pi \left(\frac{1}{\hat{R}^2} + \frac{z^2}{2\hat{R}^4} \right) \theta^J \right] + O(\hat{R}^{-6}) , \tag{3.3.29}$$

$$\frac{\delta \Delta^\mu}{\delta(\partial_0 \phi)} = \bar{\theta}^I \Gamma^\mu \left[-\frac{1}{2\hat{R}^3} y_{j'} \Gamma^{9j'} \theta^I - \frac{1}{2} \epsilon^{IJ} \Pi \left(\frac{1}{\hat{R}^2} - \frac{y^2}{2\hat{R}^4} \right) \theta^J \right] + O(\hat{R}^{-6}) . \tag{3.3.30}$$

The kinetic term in the Lagrangian (3.2.11) yields

$$\begin{aligned}
\frac{\delta \mathcal{L}_{\text{Kin}}}{\delta \dot{x}^+} &= \frac{1}{\widehat{R}^2} \left\{ p_-(x^A)^2 - \dot{x}^- + i\Delta_0^- - ip_- \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J \right\} \\
&+ \frac{1}{\widehat{R}^4} \left\{ -\frac{p_-}{2} (y^4 - z^4) + y^2 \dot{x}^- + i\Delta_2^- + i(\Delta_0^-)_{\theta^4} + \frac{i}{4} (z^2 - y^2) \Delta_0^- \right. \\
&- \frac{ip_-}{2} (z^2 - y^2) \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J - \frac{ip_-}{12} \bar{\theta}^I \Gamma^- (\mathcal{M}^2)^{IJ} \epsilon^{JL} \Pi \theta^L \\
&+ \frac{i}{4} \dot{x}^A \bar{\theta}^I \Gamma^A \left(z_k \Gamma^{-k} - y_{k'} \Gamma^{-k'} \right) \theta^I - \frac{i}{2} (\dot{x}^-) \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J + \left[-\frac{1}{2} (z^2 - y^2) \right. \\
&+ \frac{1}{2p_-^2} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] - \frac{i}{2p_-} \bar{\theta}^I \Gamma^- \partial_0 \theta^I + \frac{i}{2p_-} s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J \left. \right] \left[p_-(x^A)^2 \right. \\
&- \dot{x}^- + i\Delta_0^- - ip_- \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J \left. \right] + \left[\frac{1}{p_-^2} \dot{x}^A x'^A - \frac{i}{2p_-} \bar{\theta}^I \Gamma^- \partial_1 \theta^I \right. \\
&+ \left. \frac{i}{2p_-} s^{IJ} \bar{\theta}^I \Gamma^- \partial_0 \theta^J \right] \left(x'^- - i\Delta_0^- \right) \left. \right\} + O(\widehat{R}^{-6}) , \tag{3.3.31}
\end{aligned}$$

while the Wess-Zumino term (3.2.12) gives

$$\begin{aligned}
\frac{\delta \mathcal{L}_{\text{WZ}}}{\delta \dot{x}^+} &= \frac{i}{\widehat{R}^2} s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J + \frac{1}{\widehat{R}^4} \left\{ \frac{i}{4} s^{IJ} \bar{\theta}^I \Gamma^- (z'_j z_k \Gamma^{jk} - y'_{j'} y_{k'}) \theta^J \right. \\
&+ \frac{i}{12} s^{IJ} \bar{\theta}^I \Gamma^- (\mathcal{M}^2)^{JL} \partial_1 \theta^L - \frac{i}{4} (y^2 - z^2) s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J \\
&+ \left. \frac{i}{4} x'^A s^{IJ} \bar{\theta}^I \Gamma^A (y_{j'} \Gamma^{-j'} - z_j \Gamma^{-j}) \theta^J \right\} + O(\widehat{R}^{-6}) . \tag{3.3.32}
\end{aligned}$$

The variation is completed prior to any gauge fixing (with the worldsheet metric held fixed). After computing the variation, the lightcone coordinates x^\pm and the worldsheet metric corrections $\tilde{h}^{00}, \tilde{h}^{01}$ are to be replaced with dynamical variables according to the x^- equations of motion and the gauge conditions $x^+ = \tau$ and $T_{ab} = 0$. Hence, using a_0 and a_2 determined from the covariant gauge constraints (3.3.19, 3.3.28), we remove x^- (x^+ has already been replaced with $p_- \tau$ in the above variations) and restore proper powers of \widehat{R} in the vielbeins (so that the desired corrections enter

at $O(1/\widehat{R}^2)$). As expected, the pp-wave Hamiltonian emerges at leading order:

$$H_{\text{pp}} = \frac{p_-}{2} (x^A)^2 + \frac{1}{2p_-} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] - ip_- \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J + is^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J . \quad (3.3.33)$$

The first curvature correction to the pp-wave limit is found to be

$$\begin{aligned} H_{\text{int}} = & \frac{1}{\widehat{R}^2} \left\{ \frac{1}{4p_-} \left[y^2 (\dot{z}^2 - z'^2 - 2y'^2) + z^2 (-\dot{y}^2 + y'^2 + 2z'^2) \right] \right. \\ & + \frac{1}{8p_-^3} \left[3(\dot{x}^A)^2 - (x'^A)^2 \right] \left[(\dot{x}^A)^2 + (x'^A)^2 \right] + \frac{p_-}{8} [(x^A)^2]^2 - \frac{1}{2p_-^3} (\dot{x}^A x'^A)^2 \\ & - \frac{i}{4p_-} \sum_{a=0}^1 \bar{\theta}^I (\partial_a x^A \Gamma^A) \epsilon^{IJ} \Gamma^- \Pi (\partial_a x^B \Gamma^B) \theta^J - \frac{i}{2} p_- (x^A)^2 \bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J \\ & - \frac{i}{2p_-^2} (\dot{x}^A)^2 \bar{\theta}^I \Gamma^- \partial_0 \theta^I - \frac{ip_-}{12} \bar{\theta}^I \Gamma^- (\mathcal{M}^2)^{IJ} \epsilon^{JL} \Pi \theta^L - \frac{p_-}{2} (\bar{\theta}^I \Gamma^- \epsilon^{IJ} \Pi \theta^J)^2 \\ & - \frac{i}{2p_-^2} (\dot{x}^A x'^A) s^{IJ} \bar{\theta}^I \Gamma^- \partial_0 \theta^J - \frac{i}{4} (y^2 - z^2) s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J \\ & + \frac{i}{4} x'^A s^{IJ} \bar{\theta}^I \Gamma^A (y_{j'} \Gamma^{-j'} - z_j \Gamma^{-j}) \theta^J + \frac{i}{4} s^{IJ} \bar{\theta}^I \Gamma^- (z'_j z_k \Gamma^{jk} - y'_{j'} y_{k'} \Gamma^{j'k'}) \theta^J \\ & + \frac{i}{4p_-^2} \left[(\dot{x}^A)^2 - (x'^A)^2 \right] s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J + \frac{i}{12} s^{IJ} \bar{\theta}^I \Gamma^- (\mathcal{M}^2)^{JL} \partial_1 \theta^L \\ & \left. + \frac{1}{2} (s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J) (\bar{\theta}^K \Gamma^- \epsilon^{KL} \Pi \theta^L) + \frac{i}{4} (x^A)^2 s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J \right\} . \quad (3.3.34) \end{aligned}$$

The full Lagrangian (3.2.11, 3.2.12) can also be expressed to this order. In terms of the quantities found in equations (3.3.12, 3.3.13, 3.3.14, 3.3.23, 3.3.24), the kinetic term $\mathcal{L}_{\text{Kin}} = -\frac{1}{2} h^{ab} L_a^\mu L_b^\mu$ can be written schematically as

$$\begin{aligned} \mathcal{L}_{\text{Kin}} = & \frac{1}{2} (L_0^\mu L_0^\mu - L_1^\mu L_1^\mu)_2 + \frac{1}{2\widehat{R}^2} (L_0^\mu L_0^\mu - L_1^\mu L_1^\mu)_4 - \frac{1}{2\widehat{R}^2} \tilde{h}^{00} (L_0^\mu L_0^\mu)_2 \\ & + \frac{1}{2\widehat{R}^2} \tilde{h}^{00} (L_1^\mu L_1^\mu)_2 - \frac{1}{\widehat{R}^2} \tilde{h}^{01} (L_0^\mu L_1^\mu)_2 + O(\widehat{R}^{-4}) , \quad (3.3.35) \end{aligned}$$

where external subscripts indicate quadratic or quartic order in fields. The Wess-

Zumino term is given explicitly by:

$$\begin{aligned}
\mathcal{L}_{\text{WZ}} &= -2i\epsilon^{ab} \int_0^1 dt L_{at}^\mu s^{IJ} \bar{\theta}^I \Gamma^\mu L_{bt}^J \\
&\approx -ip_- (s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J) - \frac{i}{\widehat{R}^2} \left\{ p_- \square_2'^- + p_- (\square_0'^-)_{\theta^4} + \frac{p_-}{4} (z^2 - y^2) \square_0'^- \right. \\
&\quad \left. + \frac{1}{2} \dot{x}^- \square_0'^- - \frac{1}{2} x'^- \square_0^- + \dot{x}^A \square_1'^A - x'^A \square_1^A \right\} + O(\widehat{R}^{-4}) . \tag{3.3.36}
\end{aligned}$$

It will be useful to recast both the Hamiltonian and Lagrangian in 16-component notation (details may be found in Appendix A):

$$\begin{aligned}
H &= \frac{1}{2p_-} \left((\dot{x}^A)^2 + (x'^A)^2 + p_-^2 (x^A)^2 \right) - p_- \psi^\dagger \Pi \psi + \frac{i}{2} (\psi \psi' + \psi^\dagger \psi'^\dagger) \\
&+ \frac{1}{\widehat{R}^2} \left\{ \frac{z^2}{4p_-} [y'^2 + 2z'^2 - \dot{y}^2] - \frac{y^2}{4p_-} [z'^2 + 2y'^2 - \dot{z}^2] - \frac{1}{2p_-^3} (\dot{x}^A x'^A)^2 \right. \\
&+ \frac{1}{8p_-^3} [3(\dot{x}^A)^2 - (x'^A)^2] [(\dot{x}^A)^2 + (x'^A)^2] + \frac{p_-}{8} [(x^A)^2]^2 \\
&+ \frac{i}{8} \psi \left(z_k z'_j \gamma^{jk} - y_{k'} y'_{j'} \gamma^{j'k'} + x'^A (z_k \bar{\gamma}^A \gamma^k - y_{k'} \bar{\gamma}^A \gamma^{k'}) \right) \psi \\
&+ \frac{i}{8} \psi^\dagger \left(z_k z'_j \gamma^{jk} - y_{k'} y'_{j'} \gamma^{j'k'} + x'^A (z_k \bar{\gamma}^A \gamma^k - y_{k'} \bar{\gamma}^A \gamma^{k'}) \right) \psi^\dagger \\
&+ \frac{1}{2p_-} \left(\dot{z}^i \dot{y}^{j'} + z'^i y'^{j'} \right) \psi^\dagger \gamma^{ij'} \Pi \psi + \frac{i}{8} (z^2 - y^2) (\psi \psi' + \psi^\dagger \psi'^\dagger) \\
&+ \frac{i}{8} \left[\frac{1}{p_-^2} \left((\dot{x}^A)^2 - (x'^A)^2 \right) + (x^A)^2 \right] (\psi \psi' + \psi^\dagger \psi'^\dagger) \\
&- \frac{p_-}{2} (x^A)^2 (\psi^\dagger \Pi \psi) - \frac{i}{4p_-^2} (\dot{x}^A x'^A) (\psi \dot{\psi} + \psi^\dagger \dot{\psi}^\dagger) + \frac{p_-}{48} (\psi^\dagger \gamma^{jk} \psi) (\psi^\dagger \gamma^{jk} \psi) \\
&- \frac{p_-}{48} (\psi^\dagger \gamma^{j'k'} \psi) (\psi^\dagger \gamma^{j'k'} \psi) - \frac{i}{192} (\psi \gamma^{jk} \psi + \psi^\dagger \gamma^{jk} \psi^\dagger) (\psi^\dagger \gamma^{jk} \Pi \psi' - \psi \gamma^{jk} \Pi \psi'^\dagger) \\
&+ \frac{p_-}{2} (\psi^\dagger \Pi \psi) (\psi^\dagger \Pi \psi) + \frac{i}{192} (\psi \gamma^{j'k'} \psi + \psi^\dagger \gamma^{j'k'} \psi^\dagger) (\psi^\dagger \gamma^{j'k'} \Pi \psi' - \psi \gamma^{j'k'} \Pi \psi'^\dagger) \\
&- \frac{i}{4p_-^2} (\dot{x}^A)^2 [\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi}] - \frac{i}{4} (\psi \psi' + \psi^\dagger \psi'^\dagger) (\psi^\dagger \Pi \psi) \\
&- \frac{1}{4p_-} \left[(\dot{z}^2 - \dot{y}^2) + (z'^2 - y'^2) \right] \psi^\dagger \Pi \psi \left. \right\} + O(\widehat{R}^{-4}) . \tag{3.3.37}
\end{aligned}$$

One could scale the length of the worldsheet such that all p_- are absorbed into the upper limit on worldsheet integration over $d\sigma$. To organize correction terms by their

corresponding coupling strength in the gauge theory, however, we find it convenient to keep factors of p_- explicit in the above expression. The Lagrangian can be computed from (3.3.35, 3.3.36), giving

$$\begin{aligned}
\mathcal{L}_{\text{Kin}} = & p_- \dot{x}^- - \frac{1}{2} \left[p_-^2 (x^A)^2 - (\dot{x}^A)^2 + (x'^A)^2 \right] - \frac{ip_-}{2} (\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi}) - p_-^2 \psi \Pi \psi^\dagger \\
& + \frac{1}{2\widehat{R}^2} \left\{ (\dot{x}^-)^2 - 2p_- y^2 \dot{x}^- + \frac{1}{2} (\dot{z}^2 z^2 - \dot{y}^2 y^2) + \frac{p_-^2}{2} (y^4 - z^4) \right. \\
& - \frac{ip_-}{4} (\dot{z}_j z_k) (\psi \gamma^{jk} \psi^\dagger + \psi^\dagger \gamma^{jk} \psi) + \frac{ip_-}{4} (\dot{y}_{j'} y_{k'}) (\psi \gamma^{j'k'} \psi^\dagger + \psi^\dagger \gamma^{j'k'} \psi) \\
& - \frac{ip_-}{48} (\psi \gamma^{jk} \psi^\dagger) (\psi \gamma^{jk} \Pi \psi^\dagger - \psi^\dagger \gamma^{jk} \Pi \psi) \\
& + \frac{ip_-}{48} (\psi \gamma^{j'k'} \psi^\dagger) (\psi \gamma^{j'k'} \Pi \psi^\dagger - \psi^\dagger \gamma^{j'k'} \Pi \psi) \\
& + \frac{i}{2} \left[\frac{p_-}{2} (y^2 - z^2) - \dot{x}^- \right] (\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi}) - p_- [2\dot{x}^- - p_- (y^2 - z^2)] \psi \Pi \psi^\dagger \\
& - \frac{p_-^2}{24} (\psi^\dagger \gamma^{jk} \psi)^2 + \frac{p_-^2}{24} (\psi^\dagger \gamma^{j'k'} \psi)^2 + \frac{ip_-}{4} (\dot{x}^A z_j) (\psi \gamma^A \bar{\gamma}^j \psi^\dagger + \psi^\dagger \gamma^A \bar{\gamma}^j \psi) \\
& - \frac{ip_-}{4} (\dot{x}^A y_{j'}) (\psi \gamma^A \bar{\gamma}^{j'} \psi^\dagger + \psi^\dagger \gamma^A \bar{\gamma}^{j'} \psi) \\
& + \frac{1}{4} (\dot{x}^A \dot{x}^B) (\psi^\dagger \gamma^A \Pi \bar{\gamma}^B \psi - \psi \gamma^A \Pi \bar{\gamma}^B \psi^\dagger) \\
& - \frac{1}{2} (z'^2 z^2 - y'^2 y^2) - (x'^-)^2 + \frac{i}{2} x'^- (\psi \psi'^\dagger + \psi^\dagger \psi') \\
& - \frac{1}{4} (x'^A x'^B) (\psi^\dagger \gamma^A \Pi \bar{\gamma}^B \psi - \psi \gamma^A \Pi \bar{\gamma}^B \psi^\dagger) \\
& - \tilde{h}^{00} [2p_- \dot{x}^- - p_-^2 (x^A)^2 + (\dot{x}^A)^2 - (x'^A)^2 - ip_- (\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi}) - 2p_-^2 \psi \Pi \psi^\dagger] \\
& \left. - 2\tilde{h}^{01} [p_- x'^- + \dot{x}^A x'^A - \frac{ip_-}{2} (\psi \psi'^\dagger + \psi^\dagger \psi')] \right\} + O(\widehat{R}^{-4}) , \tag{3.3.38}
\end{aligned}$$

and

$$\begin{aligned}
\mathcal{L}_{\text{WZ}} = & -\frac{ip_-}{2} (\psi \psi' + \psi^\dagger \psi'^\dagger) - \frac{i}{\widehat{R}^2} \left\{ \frac{p_-}{8} (z'_j z_k) (\psi \gamma^{jk} \psi + \psi^\dagger \gamma^{jk} \psi^\dagger) \right. \\
& - \frac{p_-}{8} (y'_{j'} y_{k'}) (\psi \gamma^{j'k'} \psi + \psi^\dagger \gamma^{j'k'} \psi^\dagger) + \frac{1}{4} \left[\dot{x}^- + \frac{p_-}{2} (z^2 - y^2) \right] (\psi \psi' + \psi^\dagger \psi'^\dagger) \\
& - \frac{1}{4} (x'^-) (\psi \dot{\psi} + \psi^\dagger \dot{\psi}^\dagger) + \frac{i}{8} (x'^A \dot{x}^B + \dot{x}^A x'^B) (\psi^\dagger \gamma^A \Pi \bar{\gamma}^B \psi^\dagger - \psi \gamma^A \Pi \bar{\gamma}^B \psi) \\
& + \frac{p_-}{8} (x'^A z_j) (\psi^\dagger \gamma^A \bar{\gamma}^j \psi^\dagger + \psi \gamma^A \bar{\gamma}^j \psi) - \frac{p_-}{8} (x'^A y_{j'}) (\psi^\dagger \gamma^A \bar{\gamma}^{j'} \psi^\dagger + \psi \gamma^A \bar{\gamma}^{j'} \psi) \\
& \left. + \frac{p_-}{8} (\psi \gamma^{jk} \psi + \psi^\dagger \gamma^{jk} \psi^\dagger) (\psi \gamma^{jk} \Pi \psi'^\dagger - \psi^\dagger \gamma^{jk} \Pi \psi') \right\}
\end{aligned}$$

$$-\frac{p_-}{8}(\psi\gamma^{j'k'}\psi + \psi^\dagger\gamma^{j'k'}\psi^\dagger)(\psi\gamma^{j'k'}\Pi\psi'^\dagger - \psi^\dagger\gamma^{j'k'}\Pi\psi')\Big\} + O(\widehat{R}^{-4}) . \quad (3.3.39)$$

For later convenience, the Lagrangian is not fully gauge fixed, though we set \dot{x}^+ to p_- for simplicity and ignore any \ddot{x}^+ that arise through partial integration (since we will ultimately choose the lightcone gauge $x^+ = p_- \tau$). As noted above, sending $h^{00} \rightarrow -1 + \tilde{h}^{00}/\widehat{R}^2$ simply rewrites the function h^{00} , and does not amount to a particular gauge choice for the worldsheet metric.

3.4 Quantization

Our goal is to calculate explicit energy corrections due to the rather complicated perturbed Hamiltonian derived in the last section. To explain our strategy, we begin with a review of the pp-wave energy spectrum in the Penrose limit. This limit is obtained by keeping only the leading term in \widehat{R}^{-1} in the Hamiltonian expansion of (3.3.37) and leads to linear equations of motion for the fields. The eight bosonic transverse string coordinates obey the equation

$$\ddot{x}^A - x''^A + p_-^2 x^A = 0 . \quad (3.4.1)$$

This is solved by the usual expansion in terms of Fourier modes

$$\begin{aligned} x^A(\sigma, \tau) &= \sum_{n=-\infty}^{\infty} x_n^A(\tau) e^{-ik_n \sigma} , \\ x_n^A(\tau) &= \frac{i}{\sqrt{2\omega_n}} (a_n^A e^{-i\omega_n \tau} - a_{-n}^{A\dagger} e^{i\omega_n \tau}) , \end{aligned} \quad (3.4.2)$$

where $k_n = n$ (integer), $\omega_n = \sqrt{p_-^2 + k_n^2}$, and the raising and lowering operators obey the commutation relation $[a_m^A, a_n^{B\dagger}] = \delta_{mn} \delta^{AB}$. The bosonic piece of the pp-wave Hamiltonian takes the form

$$H_{\text{pp}}^B = \frac{1}{p_-} \sum_{n=-\infty}^{\infty} \omega_n \left(a_n^{A\dagger} a_n^A + 4 \right) . \quad (3.4.3)$$

The fermionic equations of motion are

$$(\dot{\psi}^\dagger + \psi') + ip_- \Pi \psi^\dagger = 0 , \quad (3.4.4)$$

$$(\dot{\psi} + \psi'^\dagger) - ip_- \Pi \psi = 0 , \quad (3.4.5)$$

where ψ is a 16-component complex $\text{SO}(9,1)$ Weyl spinor. As mentioned earlier, ψ is further restricted by a lightcone gauge fixing condition $\bar{\gamma}^9 \psi = \psi$ which reduces the number of spinor components to eight (details are given in Appendix A). In what follows, ψ and the various matrices acting on it should therefore be regarded as eight-dimensional. The fermionic equations of motion are solved by

$$\psi = \sum_{n=-\infty}^{\infty} \psi_n(\tau) e^{-ik_n \sigma} , \quad (3.4.6)$$

$$\psi_n(\tau) = \frac{1}{2\sqrt{p_-}} \left(A_n b_n e^{-i\omega_n \tau} + B_n b_{-n}^\dagger e^{i\omega_n \tau} \right) e^{-ik_n \sigma} , \quad (3.4.7)$$

$$\psi_n^\dagger(\tau) = \frac{1}{2\sqrt{p_-}} \left(\Pi B_n b_n e^{-i\omega_n \tau} - \Pi A_n b_{-n}^\dagger e^{i\omega_n \tau} \right) e^{-ik_n \sigma} , \quad (3.4.8)$$

where we have defined

$$A_n \equiv \frac{1}{\sqrt{\omega_n}} \left(\sqrt{\omega_n - k_n} - \sqrt{\omega_n + k_n} \Pi \right) , \quad (3.4.9)$$

$$B_n \equiv \frac{1}{\sqrt{\omega_n}} \left(\sqrt{\omega_n + k_n} + \sqrt{\omega_n - k_n} \Pi \right) . \quad (3.4.10)$$

The anticommuting mode operators b_n, b_n^\dagger carry a spinor index that takes eight values. In the gamma matrix representation described in Appendix A, the matrix Π is diagonal and assigns eigenvalues ± 1 to the mode operators. The fermionic canonical momentum is $\rho = ip_- \psi^\dagger$, which implies that the fermionic creation and annihilation operators obey the anticommutation rule $\{b_m^\alpha, b_n^{\beta\dagger}\} = \delta^{\alpha\beta} \delta_{mn}$. The fermionic piece of the pp-wave Hamiltonian can be written in terms of these operators as

$$H_{\text{pp}}^F = \frac{1}{p_-} \sum_{n=-\infty}^{\infty} \omega_n (b_n^{\alpha\dagger} b_n^\alpha - 4) . \quad (3.4.11)$$

Given our earlier conventions, it is necessary to invoke the coordinate reflection $x^\mu \rightarrow -x^\mu$ (Metsaev studied a similar operation on the pp-wave Hamiltonian in [25]). Such a transformation is, at this stage, equivalent to sending $x^A \rightarrow -x^A$, $p_- \rightarrow -p_-$, and $H \rightarrow -H$. In essence, this operation allows us to choose the positive-energy solutions to the fermionic equations of motion while maintaining our convention that $b^{\alpha\dagger}$ represent a creation operator and b^α denote an annihilation operator. The total pp-wave Hamiltonian

$$H_{\text{pp}} = \frac{1}{p_-} \sum_{n=-\infty}^{\infty} \omega_n \left(a_n^{A\dagger} a_n^A + b_n^{\alpha\dagger} b_n^\alpha \right) \quad (3.4.12)$$

is just a collection of free, equal mass fermionic and bosonic oscillators.

Canonical quantization requires that we express the Hamiltonian in terms of physical variables and conjugate momenta. At leading order in $1/\widehat{R}^2$, \dot{x}^A is canonically conjugate to x^A and can be expanded in terms of creation and annihilation operators. Beyond leading order, however, the conjugate variable $p_A = \delta\mathcal{L}/\delta\dot{x}^A$ differs from \dot{x}^A by terms of $O(1/\widehat{R}^2)$. Substituting these $O(1/\widehat{R}^2)$ corrected expressions for canonical momenta into the pp-wave Hamiltonian

$$H_{\text{pp}} \sim (\dot{x}^A)^2 + \psi^\dagger \Pi \psi + \psi^\dagger \psi'^\dagger \quad (3.4.13)$$

to express it as a function of canonical variables will yield indirect $O(1/\widehat{R}^2)$ corrections to the Hamiltonian (to which we must add the contribution of explicit $O(1/\widehat{R}^2)$ corrections to the action). For example, bosonic momenta in the $SO(4)$ descending from the AdS_5 subspace acquire the following corrections:

$$\begin{aligned} p_k = & \dot{z}_k + \frac{1}{\widehat{R}^2} \left\{ \frac{1}{2} y^2 p_k + \frac{1}{2p_-} \left[(p_A)^2 + (x'^A)^2 \right] p_k - \frac{1}{p_-^2} (p_A x'^A) z'_k - \frac{i}{2p_-} p_k \bar{\theta}^I \Gamma^- \partial_0 \theta^I \right. \\ & + \frac{i}{2p_-} p_k s^{IJ} \bar{\theta}^I \Gamma^- \partial_1 \theta^J - \frac{i p_-}{4} \bar{\theta}^I \Gamma^- z_j \Gamma_k^j \theta^I - \frac{i p_-}{4} \bar{\theta}^I \Gamma^k \left(z_j \Gamma^{-j} - y_{j'} \Gamma^{-j'} \right) \theta^I \\ & \left. + \frac{i}{4} p_A \epsilon^{IJ} \bar{\theta}^I \Gamma^- \left(\Gamma_k \Pi \Gamma^A + \Gamma^A \Pi \Gamma_k \right) \theta^J + \frac{i}{2p_-} z'_k \bar{\theta}^I \Gamma^- \partial_1 \theta^I - \frac{i}{2p_-} z'_k s^{IJ} \bar{\theta}^I \Gamma^- \partial_0 \theta^J \right\} \end{aligned}$$

$$+\frac{i}{4}x'^A s^{IJ} \epsilon^{JK} \bar{\theta}^I \Gamma^- \left(\Gamma_k \Pi \Gamma^A - \Gamma^A \Pi \Gamma_k \right) \theta^K \Big\} + O(\widehat{R}^{-4}) . \quad (3.4.14)$$

The leading-order relationship $p_k = \dot{z}_k$ has been substituted into the correction term at $O(1/\widehat{R}^2)$, and the lightcone gauge choice $x^+ = p_- \tau$ has been fixed after the variation.

To compute fermionic momenta $\rho = \delta \mathcal{L} / \delta \dot{\psi}$, it is convenient to work with complex 16-component spinors. Terms in \mathcal{L} relevant to the fermionic momenta ρ are as follows:

$$\begin{aligned} \mathcal{L} \sim & -ip_- \left(\psi^\dagger \dot{\psi} \right) - \frac{i}{\widehat{R}^2} \left\{ \frac{1}{4} \left[\dot{x}^- + \frac{p_-}{2} (z^2 - y^2) \right] \left(\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi} \right) \right. \\ & - \frac{p_- \tilde{h}^{00}}{2} \left(\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi} \right) + \frac{p_-}{96} \left(\psi \gamma^{jk} \dot{\psi}^\dagger \right) \left(\psi \gamma^{jk} \Pi \dot{\psi}^\dagger - \psi^\dagger \gamma^{jk} \Pi \dot{\psi} \right) \\ & \left. - \frac{x'^-}{4} \left(\psi \dot{\psi} + \psi^\dagger \dot{\psi}^\dagger \right) - (j, k \Rightarrow j', k') \right\} + O(\widehat{R}^{-4}) . \end{aligned} \quad (3.4.15)$$

This structure can be manipulated to simplify the subsequent calculation. Using partial integration, we can make the following replacement at leading order:

$$\frac{ip_-}{2} \left(\psi^\dagger \dot{\psi} + \psi \dot{\psi}^\dagger \right) = ip_- \left(\psi^\dagger \dot{\psi} \right) + \text{surface terms} . \quad (3.4.16)$$

Operations of this sort have no effect on the x^- equation of motion or the preceding calculation of $\delta \mathcal{L} / \delta \dot{x}^+$, for example. Similarly, terms in \mathcal{L} containing the matrix $(\mathcal{M})^2$ may be transformed according to

$$- \frac{ip_-}{96} \left(\psi \gamma^{jk} \dot{\psi}^\dagger \right) \left(\psi \gamma^{jk} \Pi \dot{\psi}^\dagger - \psi^\dagger \gamma^{jk} \Pi \dot{\psi} \right) = \frac{ip_-}{48} \left(\psi \gamma^{jk} \dot{\psi}^\dagger \right) \left(\psi^\dagger \gamma^{jk} \Pi \dot{\psi} \right) . \quad (3.4.17)$$

Terms of the form

$$\frac{1}{4} (\dot{x}^-) \left(\psi \dot{\psi}^\dagger + \psi^\dagger \dot{\psi} \right) , \quad (3.4.18)$$

however, cannot be treated in the same manner. The presence of (3.4.18) ultimately imposes a set of second-class constraints on the theory, and we will eventually be led

to treat ψ^\dagger as a constrained, dynamical degree of freedom in the Lagrangian. The fermionic momenta therefore take the form

$$\begin{aligned} \rho_\alpha &= ip_- \psi_\alpha^\dagger + \frac{1}{\widehat{R}^2} \left\{ \frac{i}{4} \left(\dot{x}^- + \frac{p_-}{2} (z^2 - y^2) \right) \psi_\alpha^\dagger - \frac{ip_-}{2} \tilde{h}^{00} \psi_\alpha^\dagger - \frac{ix'^-}{4} \psi_\alpha \right. \\ &\quad \left. - \frac{ip_-}{48} [(\psi \gamma^{jk} \psi^\dagger) (\psi^\dagger \gamma^{jk} \Pi)_\alpha - (j, k \rightleftharpoons j', k')] \right\} + O(\widehat{R}^{-4}) , \end{aligned} \quad (3.4.19)$$

$$\begin{aligned} \rho_\alpha^\dagger &= \frac{1}{\widehat{R}^2} \left\{ \frac{i}{4} \left(\dot{x}^- + \frac{p_-}{2} (z^2 - y^2) \right) \psi_\alpha - \frac{ip_-}{2} \tilde{h}^{00} \psi_\alpha - \frac{ix'^-}{4} \psi_\alpha^\dagger \right\} + O(\widehat{R}^{-4}) . \end{aligned} \quad (3.4.20)$$

Using (3.3.19) and (3.3.20) to replace \dot{x}^- and x'^- at leading order (in 16-component spinor notation), and using (3.3.23) to implement the appropriate curvature corrections to the h^{00} component of the worldsheet metric, we find

$$\begin{aligned} \rho &= ip_- \psi^\dagger + \frac{1}{\widehat{R}^2} \left\{ \frac{1}{4} y^2 \rho + \frac{1}{8p_-^2} [(p_A^2) + (x'^A)^2] \rho + \frac{i}{4p_-} (p_A x'^A) \psi + \frac{i}{4p_-} (\rho \Pi \psi) \rho \right. \\ &\quad \left. - \frac{i}{8p_-} (\psi \rho' + \rho \psi') \psi + \frac{i}{8p_-} \left(\psi \psi' - \frac{1}{p_-^2} \rho \rho' \right) \rho \right. \\ &\quad \left. + \frac{i}{48p_-} [(\psi \gamma^{jk} \rho) (\rho \gamma^{jk} \Pi) - (j, k, \rightleftharpoons j', k')] \right\} + O(\widehat{R}^{-4}) , \end{aligned} \quad (3.4.21)$$

$$\begin{aligned} \rho^\dagger &= \frac{1}{\widehat{R}^2} \left\{ \frac{i}{4} p_- y^2 \psi + \frac{i}{8p_-} [(p_A^2) + (x'^A)^2] \psi + \frac{1}{4p_-^2} (p_A x'^A) \rho - \frac{1}{4} (\rho \Pi \psi) \psi \right. \\ &\quad \left. - \frac{1}{8p_-^2} (\psi \rho' + \rho \psi') \rho - \frac{1}{8} \left(\psi \psi' - \frac{1}{p_-^2} \rho \rho' \right) \psi \right\} + O(\widehat{R}^{-4}) . \end{aligned} \quad (3.4.22)$$

Denoting the $O(1/\widehat{R}^2)$ corrections to ρ in (3.4.21) by Φ , the pp-wave Hamiltonian can be expressed in terms of canonical variables as

$$\begin{aligned} H_{\text{pp}} &= -p_- \psi^\dagger \Pi \psi + \frac{i}{2} \psi \psi' + \frac{i}{2} \psi^\dagger \psi'^\dagger \\ &= i \rho \Pi \psi + \frac{i}{2} \psi \psi' - \frac{i}{2p_-^2} \rho \rho' + \frac{1}{\widehat{R}^2} \left\{ \frac{i}{2p_-^2} \rho \Phi' + \frac{i}{2p_-^2} \Phi \rho' - i \Phi \Pi \psi \right\} . \end{aligned} \quad (3.4.23)$$

The $O(1/\widehat{R}^2)$ correction to the Hamiltonian can also be expressed in terms of canonical variables. The overall canonical Hamiltonian can conveniently be broken into its BMN

limit (H_{pp}), pure bosonic (H_{BB}), pure fermionic (H_{FF}) and boson-fermion (H_{BF}) interacting subsectors:

$$H_{\text{pp}} = \frac{p_-}{2}(x^A)^2 + \frac{1}{2p_-} \left[(p_A)^2 + (x'^A)^2 \right] + i\rho\Pi\psi + \frac{i}{2}\psi\psi' - \frac{i}{2p_-^2}\rho\rho' , \quad (3.4.24)$$

$$H_{\text{BB}} = \frac{1}{\widehat{R}^2} \left\{ \frac{1}{4p_-} \left[-y^2 (p_z^2 + z'^2 + 2y'^2) + z^2 (p_y^2 + y'^2 + 2z'^2) \right] + \frac{p_-}{8} [(x^A)^2]^2 \right. \\ \left. - \frac{1}{8p_-^3} \left\{ [(p_A)^2]^2 + 2(p_A)^2(x'^A)^2 + [(x'^A)^2]^2 \right\} + \frac{1}{2p_-^3} (x'^A p_A)^2 \right\} , \quad (3.4.25)$$

$$H_{\text{FF}} = -\frac{1}{4\widehat{R}^2} \left\{ \frac{1}{p_-} (\rho\Pi\psi)^2 + \frac{1}{p_-^3} (\rho\Pi\psi) \rho\rho' + \frac{1}{2p_-^3} \left(\psi\psi' - \frac{1}{p_-^2} \rho\rho' \right) \rho\rho' \right. \\ + \frac{1}{2p_-} \left(\psi\psi' - \frac{1}{p_-^2} \rho\rho' \right) (\rho\Pi\psi) + \frac{1}{2p_-^3} (\psi\rho' + \rho\psi') \rho'\psi \\ + \frac{1}{12p_-^3} (\psi\gamma^{jk}\rho) (\rho\gamma^{jk}\Pi\rho') \\ \left. - \frac{1}{48p_-} \left(\psi\gamma^{jk}\psi - \frac{1}{p_-^2} \rho\gamma^{jk}\rho \right) (\rho'\gamma^{jk}\Pi\psi - \rho\gamma^{jk}\Pi\psi') - (j, k \rightleftharpoons j', k') \right\} , \quad (3.4.26)$$

$$H_{\text{BF}} = \frac{1}{\widehat{R}^2} \left\{ \frac{i}{4} z^2 \psi\psi' - \frac{i}{8p_-^2} \left[(p_A)^2 + (x'^A)^2 \right] \psi\psi' \right. \\ + \frac{i}{4p_-^4} \left[(p_A)^2 + (x'^A)^2 + p_-^2 (y^2 - z^2) \right] \rho\rho' \\ - \frac{i}{2p_-^2} \left(p_k^2 + y'^2 - p_-^2 z^2 - \frac{1}{4} (p_A)^2 - \frac{1}{4} (x'^A)^2 - \frac{p_-^2}{2} y^2 \right) \rho\Pi\psi \\ + \frac{i}{4} (z'_j z_k) \left(\psi\gamma^{jk}\psi - \frac{1}{p_-^2} \rho\gamma^{jk}\rho \right) - \frac{i}{4} (y'_{j'} y_{k'}) \left(\psi\gamma^{j'k'}\psi - \frac{1}{p_-^2} \rho\gamma^{j'k'}\rho \right) \\ - \frac{i}{8} (z'_k y_{k'} + z_k y'_{k'}) \left(\psi\gamma^{kk'}\psi - \frac{1}{p_-^2} \rho\gamma^{kk'}\rho \right) + \frac{1}{4p_-} (p_k y_{k'} + z_k p_{k'}) \psi\gamma^{kk'}\rho \\ + \frac{1}{4p_-} (p_j z'_k) \left(\psi\gamma^{jk}\Pi\psi + \frac{1}{p_-^2} \rho\gamma^{jk}\Pi\rho \right) \\ - \frac{1}{4p_-} (p_{j'} y'_{k'}) \left(\psi\gamma^{j'k'}\Pi\psi + \frac{1}{p_-^2} \rho\gamma^{j'k'}\Pi\rho \right) \\ - \frac{1}{4p_-} (p_k y'_{k'} + z'_k p_{k'}) \left(\psi\gamma^{kk'}\Pi\psi + \frac{1}{p_-^2} \rho\gamma^{kk'}\Pi\rho \right) \\ \left. - \frac{1}{4p_-^3} (p_A x'^A) (\rho\psi' + 2\psi\rho') - \frac{i}{2p_-^2} (p_k p_{k'} - z'_k y'_{k'}) \psi\gamma^{kk'}\Pi\rho \right\} . \quad (3.4.27)$$

This Hamiltonian has one problem that we must resolve before attempting to extract its detailed consequences. At the end of Section 3.1, we argued that when the theory is restricted to the subspace of string zero-modes (i.e., excitations of the string that are independent of the worldsheet coordinate σ), curvature corrections to the leading pp-wave Hamiltonian should vanish. The only terms in the Hamiltonian that survive in this limit are those with no worldsheet spatial derivatives. Although H_{BB} has no such terms, the fermionic pieces of the Hamiltonian do. For example, H_{FF} contains a term $\widehat{R}^{-2}(\rho\Pi\psi)^2$ that would appear to modify the zero-mode spectrum at $O(1/\widehat{R}^2)$, contrary to expectation. In the end, we found that this problem can be traced to the presence of second-class constraints involving ψ^\dagger . As it turns out, the constrained quantization procedure needed to handle second-class constraints has the effect, among many others, of resolving the zero-mode paradox just outlined. To see this, we must work out the appropriate constrained quantization procedure.

The set of constraints that define canonical momenta are known as primary constraints, and take the generic form $\chi = 0$. Primary constraints can be categorized as either first or second class. Second-class constraints arise when canonical momenta do not have vanishing Poisson brackets with the primary constraints themselves: $\{\rho_\psi, \chi_\psi\} \neq 0$, $\{\rho_{\psi^\dagger}, \chi_{\psi^\dagger}\} \neq 0$. (First-class constraints are characterized by the more typical condition $\{\rho_{\psi^\dagger}, \chi_{\psi^\dagger}\} = \{\rho_\psi, \chi_\psi\} = 0$.) To the order of interest, the primary constraint equations are

$$\begin{aligned} \chi_\alpha^1 = 0 = & \rho_\alpha - ip_- \psi_\alpha^\dagger \\ & - \frac{ip_-}{8\widehat{R}^2} \left[2y^2 + \frac{1}{p_-^2} \left[(p_A)^2 + (x'^A)^2 \right] - 2(\psi^\dagger \Pi \psi) + \frac{i}{p_-} (\psi \psi' + \psi^\dagger \psi'^\dagger) \right] \psi_\alpha^\dagger \\ & - \frac{i}{4p_- \widehat{R}^2} \left[(p_A x'^A) - \frac{ip_-}{2} (\psi \psi'^\dagger + \psi^\dagger \psi') \right] \psi_\alpha + \frac{ip_-}{48\widehat{R}^2} (\psi \gamma^{jk} \psi^\dagger) (\psi^\dagger \gamma^{jk} \Pi)_\alpha, \end{aligned} \quad (3.4.28)$$

$$\chi_\alpha^2 = 0 = \rho_\alpha^\dagger - \frac{i}{4p_- \widehat{R}^2} \left[(p_A x'^A) - \frac{ip_-}{2} (\psi \psi'^\dagger + \psi^\dagger \psi') \right] \psi_\alpha^\dagger$$

$$-\frac{ip_-}{4\widehat{R}^2} \left[y^2 + \frac{1}{2p_-^2} \left[(p_A)^2 + (x'^A)^2 \right] - (\psi^\dagger \Pi \psi) + \frac{i}{2p_-} (\psi \psi' + \psi^\dagger \psi'^\dagger) \right] \psi_\alpha . \quad (3.4.29)$$

It is clear that these constraints are second-class. In the presence of second-class constraints, consistent quantization requires that the quantum anticommutator of two fermionic fields be identified with their Dirac bracket (which depends on the Poisson bracket algebra of the constraints) rather than with their classical Poisson bracket. The Dirac bracket is given in terms of Poisson brackets by (see, for example, [88])

$$\{A, B\}_D = \{A, B\}_P - \{A, \chi_N\}_P (C^{-1})^{NM} \{\chi_M, B\}_P , \quad (3.4.30)$$

where

$$C_{NM} \equiv \{\chi_N, \chi_M\}_P . \quad (3.4.31)$$

The indices N and M denote both the spinor index α and the constraint label $a = 1, 2$. For Grassmanian fields A and B , the Poisson bracket is defined by

$$\{A, B\}_P = - \left(\frac{\partial A}{\partial \psi^\alpha} \frac{\partial B}{\partial \rho_\alpha} + \frac{\partial B}{\partial \psi^\alpha} \frac{\partial A}{\partial \rho_\alpha} \right) - \left(\frac{\partial A}{\partial \psi^{\dagger\alpha}} \frac{\partial B}{\partial \rho_\alpha^\dagger} + \frac{\partial B}{\partial \psi^{\dagger\alpha}} \frac{\partial A}{\partial \rho_\alpha^\dagger} \right) . \quad (3.4.32)$$

As an example, the Dirac bracket $\{\rho_\alpha, \rho_\beta\}_D$ is readily computed (to the order of interest) by noting that the partial integration in (3.4.16) introduces an asymmetry between ψ and ψ^\dagger into the system. Since $\{\rho_\alpha, \rho_\beta\}_D$ contains

$$\{\rho_\alpha, \chi_{a\gamma}\} = O(\widehat{R}^{-2}) , \quad \{\chi_{b\eta}, \rho_\beta\} = O(\widehat{R}^{-2}) , \quad (3.4.33)$$

an immediate consequence of this asymmetry is that $\{\rho_\alpha, \rho_\beta\}_D$ vanishes to $O(1/\widehat{R}^4)$. To compute $\{\rho_\alpha, \psi_\beta\}_D$, we note that

$$\{\rho_\alpha, \chi_{(2\gamma)}\}_P = -\delta_{\alpha\rho} \frac{\partial \chi_{(2\gamma)}}{\partial \psi_\rho} = O(\widehat{R}^{-2}) , \quad (3.4.34)$$

and, to leading order,

$$(C^{-1})^{(2\gamma)(1\eta)} = -\frac{i}{p_-} \delta_{\gamma\eta} + O(\widehat{R}^{-2}) , \quad (3.4.35)$$

such that

$$\{\rho_\alpha, \psi_\beta\}_D = -\delta_{\alpha\beta} - \frac{i}{p_-} \{\rho_\alpha, \chi_{(2\beta)}\}_P . \quad (3.4.36)$$

Similar manipulations are required for $\{\psi_\alpha, \psi_\beta\}_D$, which does exhibit $O(1/\widehat{R}^2)$ corrections. The second-class constraints on the fermionic sector of the system are removed by enforcing

$$\begin{aligned} \{\rho_\alpha(\sigma), \psi_\beta(\sigma')\}_D &= -\delta_{\alpha\beta} \delta(\sigma - \sigma') + \frac{1}{4\widehat{R}^2} \delta(\sigma - \sigma') \left\{ \frac{-i}{p_-} (\rho\Pi)_\alpha \psi_\beta + \frac{i}{p_-} (\rho\Pi\psi) \delta_{\alpha\beta} \right. \\ &\quad + \frac{i}{2p_-} \left[\left(\psi\psi' \delta_{\alpha\beta} - \frac{1}{p_-} \rho\rho' \delta_{\alpha\beta} \right) + \psi'_\alpha \psi_\beta + \frac{1}{p_-^2} \rho'_\alpha \rho_\beta \right] \\ &\quad \left. + \frac{1}{2p_-^2} \left[(p_A)^2 + (x'^A)^2 \right] \delta_{\alpha\beta} + y^2 \delta_{\alpha\beta} \right\} \\ &\quad - \frac{i}{8p_- \widehat{R}^2} \left(\psi_\alpha \psi_\beta + \frac{1}{p_-^2} \rho_\alpha \rho_\beta \right) \frac{\partial}{\partial \sigma'} \delta(\sigma - \sigma') + O(\widehat{R}^{-4}) , \end{aligned} \quad (3.4.37)$$

$$\begin{aligned} \{\psi_\alpha(\sigma), \psi_\beta(\sigma')\}_D &= \frac{i}{4p_- \widehat{R}^2} \delta(\sigma - \sigma') \left\{ (\psi\Pi)_{(\alpha} \psi_{\beta)} - \frac{1}{p_-^2} (p_A x'^A) \delta_{(\alpha\beta)} \right. \\ &\quad \left. + \frac{1}{2p_-^2} [\psi'_{(\alpha} \rho_{\beta)} - \rho'_{(\alpha} \psi_{\beta)} + (\psi\rho' + \rho\psi') \delta_{(\alpha\beta)}] \right\} \\ &\quad + \frac{i}{8p_-^3 \widehat{R}^2} (\rho_{(\alpha} \psi_{\beta)} - \psi_{(\alpha} \rho_{\beta)}) \frac{\partial}{\partial \sigma'} \delta(\sigma - \sigma') + O(\widehat{R}^{-4}) , \end{aligned} \quad (3.4.38)$$

$$\{\rho_\alpha(\sigma), \rho_\beta(\sigma')\}_D = O(\widehat{R}^{-4}) . \quad (3.4.39)$$

Identifying these Dirac brackets with the quantum anticommutators of the fermionic fields in the theory naturally leads to additional $O(1/\widehat{R}^2)$ corrections to the energy spectrum. One way to implement these corrections is to retain the Fourier expansion of ψ and ψ^\dagger given in (3.4.7, 3.4.8) while transforming the fermionic creation and

annihilation operators

$$b_n^\alpha \rightarrow c_n^\alpha, \quad b_n^{\dagger\alpha} \rightarrow c_n^{\dagger\alpha}, \quad (3.4.40)$$

such that $\{\rho(c, c^\dagger), \psi(c, c^\dagger)\}_P$, for example, satisfies (3.4.37). This approach amounts to finding $O(1/\widehat{R}^2)$ corrections to $\{c_n^\alpha, c_m^{\dagger\beta}\}$ that allow the usual anticommutators to be identified with the above Dirac brackets (3.4.37-3.4.39). In practice, extracting these solutions from (3.4.37-3.4.39) can be circumvented by invoking a non-linear field redefinition $\psi \rightarrow \tilde{\psi}$, $\rho \rightarrow \tilde{\rho}$, such that

$$\{\rho(c, c^\dagger), \psi(c, c^\dagger)\}_P = \{\tilde{\rho}(b, b^\dagger), \tilde{\psi}(b, b^\dagger)\}_P. \quad (3.4.41)$$

Both representations satisfy (3.4.37), and the operators $b_n^\alpha, b_m^{\dagger\beta}$ are understood to obey the usual relations:

$$\{b_n^\alpha, b_m^{\dagger\beta}\} = \delta^{\alpha\beta} \delta_{nm}. \quad (3.4.42)$$

In general, the non-linear field redefinition $\tilde{\psi}(b, b^\dagger) = \psi(b, b^\dagger) + \dots$ contains corrections that are cubic in the fields $\rho(b, b^\dagger)$, $\psi(b, b^\dagger)$, $x^A(a, a^\dagger)$ and $p_A(a, a^\dagger)$. Such correction terms can be written down by inspection, with matrix-valued coefficients to be solved for by comparing $\{\tilde{\rho}(b, b^\dagger), \tilde{\psi}(b, b^\dagger)\}_P$ and $\{\tilde{\psi}(b, b^\dagger), \tilde{\psi}(b, b^\dagger)\}_P$ with (3.4.37, 3.4.38). A straightforward computation yields

$$\begin{aligned} \rho_\alpha \rightarrow \tilde{\rho}_\alpha &= \rho_\alpha, & (3.4.43) \\ \psi_\beta \rightarrow \tilde{\psi}_\beta &= \psi_\beta + \frac{i}{8p_- \widehat{R}^2} \left\{ (\psi' \psi) \psi_\beta - 2(\rho \Pi \psi) \psi_\beta - \frac{1}{p_-^2} (\rho' \rho) \psi_\beta + \frac{2}{p_-^2} (p_A x'^A) \rho_\beta \right. \\ &\quad \left. + \frac{1}{p_-^2} [(\rho' \psi) \rho_\beta - (\rho \psi') \rho_\beta] + 2ip_- \left[y^2 \psi_\beta + \frac{1}{2p_-^2} \left((p_A)^2 + (x'^A)^2 \right) \psi_\beta \right] \right\}. & (3.4.44) \end{aligned}$$

This approach to enforcing the modified Dirac bracket structure amounts to adding

$O(1/\widehat{R}^2)$ correction terms to the Hamiltonian while keeping the standard commutation relations. It is much more convenient for calculating matrix elements than the alternative approach of adding $O(1/\widehat{R}^2)$ operator corrections to the fermi field anticommutators $\{b, b^\dagger\}$.

By invoking the redefinitions in (3.4.43, 3.4.44), the pieces of the interaction Hamiltonian that involve fermions take the final forms

$$\begin{aligned}
H_{\text{FF}} = & -\frac{1}{4p_-^3 \widehat{R}^2} \left\{ p_-^2 \left[(\psi' \psi) + \frac{1}{p_-^2} (\rho \rho') \right] (\rho \Pi \psi) - \frac{p_-^2}{2} (\psi' \psi)^2 + (\psi' \psi) (\rho' \rho) \right. \\
& - \frac{1}{2p_-^2} (\rho' \rho)^2 + (\rho \psi') (\rho' \psi) - \frac{1}{2} [(\psi \rho') (\psi \rho') + (\psi' \rho)^2] + \frac{1}{12} (\psi \gamma^{jk} \rho) (\rho \gamma^{jk} \Pi \rho') \\
& \left. - \frac{p_-^2}{48} \left(\psi \gamma^{jk} \psi - \frac{1}{p_-^2} \rho \gamma^{jk} \rho \right) (\rho' \gamma^{jk} \Pi \psi - \rho \gamma^{jk} \Pi \psi') - (j, k \rightleftharpoons j', k') \right\}, \tag{3.4.45}
\end{aligned}$$

$$\begin{aligned}
H_{\text{BF}} = & \frac{1}{\widehat{R}^2} \left\{ -\frac{i}{4p_-^2} \left[(p_A)^2 + (x'^A)^2 + p_-^2 (y^2 - z^2) \right] \left(\psi \psi' - \frac{1}{p_-^2} \rho \rho' \right) \right. \\
& - \frac{1}{2p_-^3} (p_A x'^A) (\rho \psi' + \psi \rho') - \frac{i}{2p_-^2} \left(p_k^2 + y'^2 - p_-^2 z^2 \right) \rho \Pi \psi \\
& + \frac{i}{4} (z'_j z_k) \left(\psi \gamma^{jk} \psi - \frac{1}{p_-^2} \rho \gamma^{jk} \rho \right) - \frac{i}{4} (y'_{j'} y_{k'}) \left(\psi \gamma^{j'k'} \psi - \frac{1}{p_-^2} \rho \gamma^{j'k'} \rho \right) \\
& - \frac{i}{8} (z'_k y_{k'} + z_k y'_{k'}) \left(\psi \gamma^{kk'} \psi - \frac{1}{p_-^2} \rho \gamma^{kk'} \rho \right) + \frac{1}{4p_-} (p_k y_{k'} + z_k p_{k'}) \psi \gamma^{kk'} \rho \\
& + \frac{1}{4p_-} (p_j z'_k) \left(\psi \gamma^{jk} \Pi \psi + \frac{1}{p_-^2} \rho \gamma^{jk} \Pi \rho \right) \\
& - \frac{1}{4p_-} (p_{j'} y'_{k'}) \left(\psi \gamma^{j'k'} \Pi \psi + \frac{1}{p_-^2} \rho \gamma^{j'k'} \Pi \rho \right) \\
& - \frac{1}{4p_-} (p_k y'_{k'} + z'_k p_{k'}) \left(\psi \gamma^{kk'} \Pi \psi + \frac{1}{p_-^2} \rho \gamma^{kk'} \Pi \rho \right) \\
& \left. - \frac{i}{2p_-^2} (p_k p_{k'} - z'_k y'_{k'}) \psi \gamma^{kk'} \Pi \rho \right\}. \tag{3.4.46}
\end{aligned}$$

The full Hamiltonian is the sum of these two terms plus the bosonic interaction term H_{BB} (3.4.25) and the free Hamiltonian H_{pp} (3.4.24). This system is quantized by imposing the standard (anti)commutator algebra for x^A, ψ and their conjugate variables p^A, ρ . This will be done by expanding the field variables in creation and

annihilation operators in a standard way.

Returning to the phenomenon that led us to explore second-class constraints in the first place, note that (3.4.45) manifestly vanishes on the subspace of string zero-modes because all terms have at least one worldsheet spatial derivative. The bose-fermi mixing Hamiltonian (3.4.46) still has terms that can lead to curvature corrections to the string zero-mode energies, but their net effect vanishes by virtue of nontrivial cancellations between terms that split $SO(4) \times SO(4)$ indices and terms that span the entire $SO(8)$. How this comes about will be seen when we actually compute matrix elements of this Hamiltonian.

3.5 Energy spectrum

To compute the energy spectrum correct to first order in $O(\hat{R}^{-1})$, we will do degenerate first-order perturbation theory on the Fock space of eigenstates of the free Hamiltonian H_{pp} . The degenerate subspaces of the BMN theory are spanned by fixed numbers of creation operators with specified mode indices (subject to the constraint that the mode indices sum up to zero) acting on the ground state $|J\rangle$, where $J = p_- \hat{R}^2$ is the angular momentum (assumed large) of the string center of mass in its motion around the equator of the S^5 . In this chapter we restrict attention to “two-impurity states” generated by pairs of creation operators of equal and opposite mode number. For each positive mode number n , the 16 bosonic and fermionic creation operators can be combined in pairs to form the following 256 degenerate two-impurity states:

$$a_n^{A\dagger} a_{-n}^{B\dagger} |J\rangle, \quad b_n^{\alpha\dagger} b_{-n}^{\beta\dagger} |J\rangle, \quad a_n^{A\dagger} b_{-n}^{\alpha\dagger} |J\rangle, \quad a_{-n}^{A\dagger} b_n^{\alpha\dagger} |J\rangle. \quad (3.5.1)$$

The creation operators are classified under the residual $SO(4) \times SO(4)$ symmetry to which the isometry group of the $AdS_5 \times S^5$ target space is broken by the lightcone gauge quantization procedure. The bosonic creation operators $a_n^{A\dagger}$ decompose as $(\mathbf{4}, \mathbf{1}) + (\mathbf{1}, \mathbf{4})$, or, in the $SU(2)^2 \times SU(2)^2$ notation introduced in [26],

as $(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$. Analogously, the fermionic operators $b_n^{\alpha\dagger}$ decompose as $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) + (\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ under the covering group. It is useful to note that the two fermion irreps are eigenvectors, with opposite eigenvalue, of the Π operator introduced in (3.2.6). To find the perturbed energy spectrum, we must compute explicit matrix elements of H_{int} in this basis and then diagonalize the resulting 256×256 matrix. We will compare the perturbed energy eigenvalues with general expectations from $PSU(2, 2|4)$ as well as with the large R -charge limit of the anomalous dimensions of gauge theory operators with two R -charge defects. Higher-impurity string states can be treated in the same way, but we defer such questions to a later chapter. Our purpose here is primarily to check that our methods (choice of action, lightcone gauge reduction, quantization rules, etc.) are consistent and correct. Due to the algebraic complexity met with at each step, this check is far from trivial. Once reassured on these fundamental points, we can go on to examine a wider range of physically interesting issues.

The first step in carrying out this program is to expand H_{int} in creation and annihilation operators using (3.4.2, 3.4.7) for x^A, ψ and the related expansions for p^A, ρ . As an example, we quote the result for H_{BB} (keeping only terms with two creation and two annihilation operators):

$$\begin{aligned}
H_{\text{BB}} = & -\frac{1}{32p_- \widehat{R}^2} \sum \frac{\delta(n+m+l+p)}{\xi} \times \\
& \left\{ 2 \left[\xi^2 - (p_-^4 - k_l k_p k_n k_m) + \omega_n \omega_m k_l k_p + \omega_l \omega_p k_n k_m + 2\omega_n \omega_l k_m k_p \right. \right. \\
& + 2\omega_m \omega_p k_n k_l \left. \right] a_{-n}^{\dagger A} a_{-m}^{\dagger A} a_l^B a_p^B + 4 \left[\xi^2 - (p_-^4 - k_l k_p k_n k_m) - 2\omega_n \omega_m k_l k_p \right. \\
& + \omega_l \omega_m k_n k_p - \omega_n \omega_l k_m k_p - \omega_m \omega_p k_n k_l + \omega_n \omega_p k_m k_l \left. \right] a_{-n}^{\dagger A} a_{-l}^{\dagger B} a_m^A a_p^B \\
& + 2 \left[8k_l k_p a_{-n}^{\dagger i} a_{-l}^{\dagger j} a_m^i a_p^j + 2(k_l k_p + k_n k_m) a_{-n}^{\dagger i} a_{-m}^{\dagger i} a_l^j a_p^j \right. \\
& + (\omega_l \omega_p + k_l k_p - \omega_n \omega_m - k_n k_m) a_{-n}^{\dagger i} a_{-m}^{\dagger i} a_l^{j'} a_p^{j'} - 4(\omega_l \omega_p - k_l k_p) a_{-n}^{\dagger i} a_{-l}^{\dagger j'} a_m^i a_p^{j'} \\
& \left. \left. - (i, j \rightleftharpoons i', j') \right] \right\}, \tag{3.5.2}
\end{aligned}$$

with $\xi \equiv \sqrt{\omega_n \omega_m \omega_l \omega_p}$. The expansion of the interaction terms involving fermi fields are too complicated to be worth writing down explicitly at this stage. Schematically, we organize the two-impurity matrix elements of the perturbing Hamiltonian as shown in Table 3.1.

H_{int}	$a_n^{A\dagger} a_{-n}^{B\dagger} J\rangle$	$b_n^{\alpha\dagger} b_{-n}^{\beta\dagger} J\rangle$	$a_n^{A\dagger} b_{-n}^{\alpha\dagger} J\rangle$	$a_{-n}^{A\dagger} b_n^{\alpha\dagger} J\rangle$
$\langle J a_n^A a_{-n}^B$	H_{BB}	H_{BF}	0	0
$\langle J b_n^\alpha b_{-n}^\beta$	H_{BF}	H_{FF}	0	0
$\langle J a_n^A b_{-n}^\alpha$	0	0	H_{BF}	\mathcal{H}_{BF}
$\langle J a_{-n}^A b_n^\alpha$	0	0	H_{BF}	H_{BF}

Table 3.1: Structure of the matrix of first-order energy perturbations in the space of two-impurity string states

To organize the perturbation theory, it is helpful to express everything in terms of two parameters: J and λ' . In the duality between Type IIB superstring theory on $AdS_5 \times S^5$ and $\mathcal{N} = 4$ $SU(N_c)$ super Yang-Mills theory in four dimensions, we identify

$$\begin{aligned}
\mathcal{N} = 4 \text{ SYM} & \quad AdS_5 \times S^5 , \\
SU(N_c) & \quad \Longleftrightarrow \quad \int_{S^5} F_5 = N_c , \\
g_{YM}^2 N_c & \quad \Longleftrightarrow \quad \widehat{R}^4 , \\
g_{YM}^2 & \quad \Longleftrightarrow \quad g_s .
\end{aligned} \tag{3.5.3}$$

In the pp-wave limit, however, the AdS/CFT dictionary reads

$$\begin{aligned}
R & \quad \Longleftrightarrow \quad p_- \widehat{R}^2 = J , \\
\frac{R^2}{N_c} & \quad \Longleftrightarrow \quad g_s p_-^2 = g_2 , \\
R \rightarrow \infty & \quad \Longleftrightarrow \quad p_- \widehat{R}^2, N_c \rightarrow \infty .
\end{aligned} \tag{3.5.4}$$

The modified 't Hooft coupling

$$\lambda' = \frac{g_{\text{YM}}^2 N_c}{R^2} \rightleftharpoons \frac{1}{p_-^2} \quad (3.5.5)$$

is kept fixed in the $R, N_c \rightarrow \infty$ limit. (We have kept $\alpha' = \mu = 1$.) Since the gauge theory is perturbative in $\lambda = g_{\text{YM}}^2 N_c$, and p_-^2 on the string side is mapped to $R^2/(g_{\text{YM}}^2 N_c)$, we will expand string energies ω_q in powers of $1/p_-$, keeping terms up to some low order to correspond with the loop expansion in the gauge theory. This type of dictionary would be incorrect in the original coordinate system characterized by the lightcone coordinates $t = x^+ - (x^-/2\hat{R}^2)$ and $\phi = x^+ + (x^-/2\hat{R}^2)$ given in (3.1.4). In this case, one would calculate corrections to $R \rightleftharpoons p_- \hat{R}^2$ appearing in the perturbing Hamiltonian (which amount to operator-valued corrections to p_-).

3.5.1 Evaluating Fock space matrix elements of H_{BB}

We now proceed to the construction of the perturbing Hamiltonian matrix on the space of degenerate two-impurity states. To convey a sense of what is involved, we display the matrix elements of H_{BB} (3.4.25) between the bosonic two-impurity Fock space states:

$$\begin{aligned} \left\langle J \left| a_n^A a_{-n}^B (H_{\text{BB}}) a_{-n}^{C\dagger} a_n^{D\dagger} \right| J \right\rangle &= (N_{\text{BB}}(n^2 \lambda') - 2n^2 \lambda') \frac{\delta^{AD} \delta^{BC}}{J} \\ &+ \frac{n^2 \lambda'}{J(1 + n^2 \lambda')} [\delta^{ab} \delta^{cd} + \delta^{ad} \delta^{bc} - \delta^{ac} \delta^{bd}] \\ &- \frac{n^2 \lambda'}{J(1 + n^2 \lambda')} [\delta^{a'b'} \delta^{c'd'} + \delta^{a'd'} \delta^{b'c'} - \delta^{a'c'} \delta^{b'd'}] \\ &\approx (n_{\text{BB}} - 2) \frac{n^2 \lambda'}{J} \delta^{AD} \delta^{BC} + \frac{n^2 \lambda'}{J} [\delta^{ab} \delta^{cd} + \delta^{ad} \delta^{bc} - \delta^{ac} \delta^{bd}] \\ &- \frac{n^2 \lambda'}{J} [\delta^{a'b'} \delta^{c'd'} + \delta^{a'd'} \delta^{b'c'} - \delta^{a'c'} \delta^{b'd'}] + O(\lambda'^2), \quad (3.5.6) \end{aligned}$$

where lower-case $SO(4)$ indices $a, b, c, d \in 1, \dots, 4$ indicate that A, B, C, D are chosen from the first $SO(4)$, and $a', b', c', d' \in 5, \dots, 8$ indicate the second $SO(4)$ ($A, B, C, D \in 5, \dots, 8$). We have also displayed the further expansion of these $O(1/J)$ matrix

elements in powers of λ' (using the basic BMN-limit energy eigenvalue condition $\omega_n/p_- = \sqrt{1 + \lambda' n^2}$). This is to facilitate eventual contact with perturbative gauge theory via AdS/CFT duality. Note that H_{BB} does not mix states built out of oscillators from different $SO(4)$ subgroups. There is a parallel no-mixing phenomenon in the gauge theory: two-impurity bosonic operators carrying spacetime vector indices do not mix with spacetime scalar bosonic operators carrying R -charge vector indices.

Due to operator ordering ambiguities, two-impurity matrix elements of H_{BB} can differ by contributions proportional to $\delta^{AD}\delta^{BC}$, depending on the particular prescription chosen [26]. $N_{\text{BB}}(n^2\lambda')$ is an arbitrary function of $n^2\lambda'$, which is included to account for such ambiguities (we will shortly succeed in fixing it). To match the dual gauge theory physics, it is best to expand N_{BB} as a power series in λ' . The zeroth-order term must vanish if the energy correction is to be perturbative in the gauge coupling. The next term in the expansion contributes one arbitrary constant (the n_{BB} term) and each higher term in the λ' expansion in principle contributes one additional arbitrary constant to this sector of the Hamiltonian. Simple general considerations will fix them all.

3.5.2 Evaluating Fock space matrix elements of H_{FF}

The calculation of the two-impurity matrix elements of the parts of H_{int} that involve fermionic fields is rather involved and we found it necessary to employ symbolic manipulation programs to keep track of the many different terms. The end results are fairly concise, however. For H_{FF} we find

$$\begin{aligned} \left\langle J \left| b_n^\alpha b_{-n}^\beta (H_{\text{FF}}) b_{-n}^{\gamma\dagger} b_n^{\delta\dagger} \right| J \right\rangle &= (N_{\text{FF}}(n^2\lambda') - 2n^2\lambda') \frac{\delta^{\alpha\delta}\delta^{\beta\gamma}}{J} \\ &+ \frac{n^2\lambda'}{24J(1+n^2\lambda')} \left[(\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\beta\gamma} + (\gamma^{ij})^{\alpha\beta} (\gamma^{ij})^{\gamma\delta} - (\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\beta\delta} \right] \\ &- \frac{n^2\lambda'}{24J(1+n^2\lambda')} \left[(\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\beta\gamma} + (\gamma^{i'j'})^{\alpha\beta} (\gamma^{i'j'})^{\gamma\delta} - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\beta\delta} \right] \end{aligned}$$

$$\begin{aligned}
\approx & (n_{\text{FF}} - 2) \frac{n^2 \lambda'}{J} \delta^{\alpha\delta} \delta^{\beta\gamma} + \frac{n^2 \lambda'}{24J} [(\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\beta\gamma} + (\gamma^{ij})^{\alpha\beta} (\gamma^{ij})^{\gamma\delta} - (\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\beta\delta}] \\
& - \frac{n^2 \lambda'}{24J} [(\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\beta\gamma} + (\gamma^{i'j'})^{\alpha\beta} (\gamma^{i'j'})^{\gamma\delta} - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\beta\delta}] + O(\lambda'^2) .
\end{aligned}
\tag{3.5.7}$$

This sector has its own normal-ordering function N_{FF} , with properties similar those of N_{BB} described above. The index structure of the fermionic matrix elements is similar to that of its bosonic counterpart (3.5.6).

We will now introduce some useful projection operators that will help us understand the selection rules implicit in the index structure of (3.5.7). The original 16-component spinors ψ were reduced to eight components by the Weyl condition $\bar{\gamma}^9 \psi = \psi$. The remaining eight components are further divided into spinors $\tilde{\psi}$ and $\hat{\psi}$, which are even or odd under the action of Π :

$$\begin{aligned}
\Pi \tilde{\psi} &= -\tilde{\psi} , & \Pi \tilde{b}^{\dagger\alpha} &= -\tilde{b}^{\dagger\alpha} , \\
\Pi \hat{\psi} &= \hat{\psi} , & \Pi \hat{b}^{\dagger\alpha} &= \hat{b}^{\dagger\alpha} .
\end{aligned}
\tag{3.5.8}$$

The spinors $\hat{\psi}$ transform in the $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ of $SO(4) \times SO(4)$, while $\tilde{\psi}$ transform in the $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$. This correlation between Π -parity and $SO(4) \times SO(4)$ representation will be very helpful for analyzing complicated fermionic matrix elements.

We denote the $SU(2)$ generators of the active factors of the $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ irrep as Σ^+ and Ω^+ , where the Σ act on the $SO(4)$ descended from the AdS_5 , and the Ω act on the $SO(4)$ coming from the S^5 . The $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ generators are similarly labeled by Σ^- and Ω^- . Each set of spinors is annihilated by its counterpart set of $SU(2)$ generators:

$$\begin{aligned}
\Sigma^+ \hat{b}^{\dagger\alpha} &= \Omega^+ \hat{b}^{\dagger\alpha} = 0 , \\
\Sigma^- \tilde{b}^{\dagger\alpha} &= \Omega^- \tilde{b}^{\dagger\alpha} = 0 .
\end{aligned}
\tag{3.5.9}$$

In terms of the projection operators

$$\Pi_+ = \frac{1}{2}(1 + \Pi) , \quad \Pi_- = \frac{1}{2}(1 - \Pi) , \quad (3.5.10)$$

which select the disjoint $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ and $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ irreps, respectively, we have

$$\begin{aligned} \Pi_+ \psi &= \hat{\psi} , & \Pi_+ \hat{b}^\alpha &= \hat{b}^\alpha , \\ \Pi_- \psi &= \tilde{\psi} , & \Pi_- \tilde{b}^\alpha &= \tilde{b}^\alpha . \end{aligned} \quad (3.5.11)$$

The Π_\pm projections commute with the $SO(4)$ generator matrices $\gamma^{ij}, \gamma^{i'j'}$, a fact that implies certain useful selection rules for the one-loop limit of (3.5.7). The rules are most succinctly stated using an obvious \pm shorthand to indicate the representation content of states created by multiple fermionic creation operators. In brief, one finds that $++$ states connect only with $++$ and $--$ states connect only with $--$. The only subtle point is the statement that all $++ \rightarrow --$ matrix elements of (3.5.7) must vanish: this is the consequence of a simple cancellation between two terms. This observation will simplify the matrix diagonalization we will eventually carry out.

3.5.3 Evaluating Fock space matrix elements of H_{BF}

The H_{BF} sector in the Hamiltonian mediates mixing between spacetime bosons of the two types (pure boson and bi-fermion) as well as between spacetime fermions (which of course contain both bosonic and fermionic oscillator excitations). The 64-dimensional boson mixing matrix

$$\left\langle J \left| b_n^\alpha b_{-n}^\beta (H_{\text{BF}}) a_{-n}^{A\dagger} a_n^{B\dagger} \right| J \right\rangle ,$$

is an off-diagonal block in the bosonic sector of the perturbation matrix in Table 3.1. The same methods used earlier in this section to reduce Fock space matrix elements involving fermi fields can be used here to obtain the simple explicit result (we omit

the details)

$$\begin{aligned}
\left\langle J \left| b_n^\alpha b_{-n}^\beta (H_{\text{BF}}) a_{-n}^{A\dagger} a_n^{B\dagger} \right| J \right\rangle &= \frac{n^2 \lambda'}{2J(1+n^2 \lambda')} \left\{ \sqrt{1+n^2 \lambda'} \left[(\gamma^{ab'})^{\alpha\beta} - (\gamma^{a'b})^{\alpha\beta} \right] \right. \\
&\quad \left. + n\sqrt{\lambda'} \left[(\gamma^{a'b'})^{\alpha\beta} - (\gamma^{ab})^{\alpha\beta} + (\delta^{ab} - \delta^{a'b'}) \delta^{\alpha\beta} \right] \right\} \\
&\approx \frac{n^2 \lambda'}{2J} \left[(\gamma^{ab'})^{\alpha\beta} - (\gamma^{a'b})^{\alpha\beta} \right] + O(\lambda'^{3/2}) . \quad (3.5.12)
\end{aligned}$$

The complex conjugate of this matrix element gives the additional off-diagonal component of the upper 128×128 block of spacetime bosons. We note that terms in the H_{BF} sector split the $SO(8)$ group (manifest in the pp-wave limit) into its $SO(4)$ constituents such that states of the form $a_{-n}^{a'\dagger} a_n^{b'\dagger} |J\rangle$, for example, which descend strictly from the S^5 subspace, vanish in this subsector. This behavior is reproduced in the gauge theory, wherein two-boson states that are either spacetime scalars or scalars of the R -charge group do not mix with bi-fermionic scalars in either irrep.

The 128-dimensional subsector of spacetime fermions is mixed by matrix elements of the same Hamiltonian taken between fermionic string states of the general form $b_n^{\alpha\dagger} a_{-n}^{A\dagger} |J\rangle$. Our standard methods yield the following simple results for the two independent types of spacetime fermion mixing matrix elements:

$$\begin{aligned}
\left\langle J \left| b_n^\alpha a_{-n}^A (H_{\text{BF}}) b_n^{\beta\dagger} a_{-n}^{B\dagger} \right| J \right\rangle &= N_{\text{BF}}(n^2 \lambda') \frac{\delta^{AB} \delta^{\alpha\beta}}{J} \\
&+ \frac{n^2 \lambda'}{2J(1+n^2 \lambda')} \left\{ (\gamma^{ab})^{\alpha\beta} - (\gamma^{a'b'})^{\alpha\beta} - (3+4n^2 \lambda') \delta^{ab} \delta^{\alpha\beta} - (5+4n^2 \lambda') \delta^{a'b'} \delta^{\alpha\beta} \right\} \\
&\approx \frac{n^2 \lambda'}{2J} \left\{ (\gamma^{ab})^{\alpha\beta} - (\gamma^{a'b'})^{\alpha\beta} + [(2n_{\text{BF}}-3)\delta^{ab} + (2n_{\text{BF}}-5)\delta^{a'b'}] \delta^{\alpha\beta} \right\} + O(\lambda'^2) , \quad (3.5.13)
\end{aligned}$$

$$\begin{aligned}
\left\langle J \left| b_n^\alpha a_{-n}^A (H_{\text{BF}}) b_{-n}^{\beta\dagger} a_n^{B\dagger} \right| J \right\rangle &= \frac{n^2 \lambda'}{2J\sqrt{1+n^2 \lambda'}} \left\{ (\gamma^{ab})^{\alpha\beta} - (\gamma^{a'b'})^{\alpha\beta} \right. \\
&\quad \left. - \frac{n\lambda'^{1/2}}{\sqrt{1+n^2 \lambda'}} \left[(\gamma^{ab'})^{\alpha\beta} - (\gamma^{a'b})^{\alpha\beta} \right] - \delta^{\alpha\beta} (\delta^{ab} - \delta^{a'b'}) \right\}
\end{aligned}$$

$$\approx \frac{n^2 \lambda'}{2J} \left\{ (\gamma^{ab})^{\alpha\beta} - (\gamma^{a'b'})^{\alpha\beta} - (\delta^{ab} - \delta^{a'b'}) \delta^{\alpha\beta} \right\} + O(\lambda'^{3/2}) . \quad (3.5.14)$$

Equation (3.5.13) involves yet another normal-ordering function. Since these functions have a nontrivial effect on the spectrum, we must give them specific values before we can calculate actual numerical eigenvalues. The key point is that the structure of the perturbing Hamiltonian implies certain relations between all the normal-ordering functions. Because the interaction Hamiltonian is quartic in oscillators, normal-ordering ambiguities give rise to terms quadratic in oscillators, appearing as constant contributions to the diagonal matrix elements. There are normal-ordering contributions from each sector of the theory: H_{BB} contributes a single term quadratic in bosonic oscillators; H_{FF} yields a term quadratic in fermionic oscillators; H_{BF} contributes one term quadratic in bosons and one quadratic in fermions. The bosonic contributions multiply terms of the form $a^\dagger a$, which are collected into the function $N_{\text{BB}}(n^2 \lambda')$ with one contribution from H_{BB} and one contribution from H_{BF} . Similarly, $N_{\text{FF}}(n^2 \lambda')$ collects terms multiplying $b^\dagger b$, receiving one contribution from H_{FF} and one contribution from H_{BF} . Normal-ordering contributions from both $a^\dagger a$ and $b^\dagger b$ terms are non-vanishing in the spacetime fermion subsector; all possible normal-ordering ambiguities appear in this subspace. The normal-ordering function $N_{\text{BF}}(n^2 \lambda')$ therefore must satisfy

$$N_{\text{BF}}(n^2 \lambda') = N_{\text{BB}}(n^2 \lambda') + N_{\text{FF}}(n^2 \lambda') . \quad (3.5.15)$$

The normal ordering functions are basically finite renormalizations that must be adjusted so that the spectrum reflects the $PSU(2, 2|4)$ global supersymmetry of the classical worldsheet action (a symmetry we want to preserve at the quantum level).

As has been explained elsewhere [26, 28] (and as we shall shortly review), energy levels should be organized into multiplets obtained by acting on a “highest-weight” level with all possible combinations of the eight R -charge raising supercharges. All the states obtained by acting with a total of L supercharges have the same energy and we

will refer to them as states at level L in the supermultiplet. The levels of a multiplet run from $L = 0$ to $L = 8$. A careful inspection of the way the normal ordering functions contribute to the energies of states in the two-impurity sector shows that states at levels $L = 0, 8$ are shifted by N_{BB} only. Similarly, levels $L = 2, 4, 6$ are shifted by N_{FF} or N_{BB} and one must have $N_{BB} = N_{FF}$ if those levels are to remain internally degenerate. Finally, levels $L = 1, 3, 5, 7$ are shifted by N_{BF} only. By supersymmetry, the level spacing must be uniform throughout the supermultiplet and this is only possible if we also set $N_{BB} = N_{BF}$. But then the constraint $N_{BF} = N_{BB} + N_{FF}$ can only be met by setting $N_{BB} = N_{FF} = N_{BF} = 0$, which then eliminates any normal-ordering ambiguity from the string theory. This is basically an exercise in using global symmetry conditions to fix otherwise undetermined finite renormalizations.

3.5.4 Diagonalizing the one-loop perturbation matrix

We are now ready to diagonalize the perturbing Hamiltonian and examine whether the resulting energy shifts have the right multiplet structure and whether the actual eigenvalues match gauge theory expectations. To simplify the problem, we will begin by diagonalizing the perturbation matrix expanded to first nontrivial order in both $1/J$ and λ' . Our results should, by duality, match one-loop gauge theory calculations and we will eventually return to the problem of finding the string spectrum correct to higher orders in λ' . From the structure of the results just obtained for the perturbation matrices, we can see that the general structure of the energy eigenvalues of two-impurity states must be

$$E_{\text{int}}(n) = 2 + n^2 \lambda' \left(1 + \frac{\Lambda}{J} + O(J^{-2}) \right) + O(\lambda'^2) , \quad (3.5.16)$$

where Λ is dimensionless and the dependence on $1/J$, λ' and mode number n is given by (3.5.6, 3.5.7). The eigenvalues Λ must meet certain conditions if the requirements of $PSU(2,2|4)$ symmetry are to be met, and we will state those conditions before solving the eigenvalue problem.

The eigenvalues Λ must meet certain conditions if the requirements of $PSU(2, 2|4)$ symmetry are to be met. The eigenvalues in question are lightcone energies and thus dual to the gauge theory quantity $\Delta = D - J$, the difference between scaling dimension and R -charge. Since conformal invariance is part of the full symmetry group, states are organized into conformal multiplets built on conformal primaries. A supermultiplet will contain several conformal primaries having the same value of Δ and transforming into each other under the supercharges. All 16 supercharges increment the dimension of an operator by $1/2$, but only eight of them (call them \mathcal{Q}_α) also increment the R -charge by $1/2$, so as to leave Δ unchanged. These eight supercharges act as “raising operators” on the conformal primaries of a supermultiplet: starting from a super-primary of lowest R -charge, the other conformal primaries are created by acting on it in all possible ways with the eight \mathcal{Q}_α . Primaries obtained by acting with L factors of \mathcal{Q}_α on the super-primary are said to be at level L in the supermultiplet (since the \mathcal{Q}_α anticommute, the range is $L = 0$ to $L = 8$). The multiplicities of states at the various levels are also determined: for each $L = 0$ primary operator, there will be C_L^8 such operators at level L (where C_m^n is the binomial coefficient). If the $L = 0$ primary has multiplicity s , summing over all L gives $2^8 s = 256s$ conformal primaries in all.

These facts severely restrict the quantity Λ in the general expression (3.5.16) above. Although the states in the degenerate multiplet all have the same J , they actually belong to different levels L in more than one supermultiplet. A state of given L is a member of a supermultiplet built on a “highest-weight” or super-primary state with $R = J - L/2$. Since all the primaries in a supermultiplet have the same Δ , the joint dependence of eigenvalues on λ, J, L must be of the form $\Delta(\lambda, J - L/2)$. The only way the expansion of (3.5.16) can be consistent with this is if $\Lambda = L + c$, where c is a pure numerical constant (recall that $\lambda' = \lambda/J^2$). Successive members of a supermultiplet must therefore have eigenvalues separated by exactly one and the difference between “top” ($L = 8$) and “bottom” ($L = 0$) eigenvalues for Λ must be exactly eight.

3.5.5 Details of the one-loop diagonalization procedure.

We now confront the problem of explicitly diagonalizing the first-order perturbation matrix Λ (obtained by expanding the relevant matrix elements to first order in λ'). The matrix block diagonalizes on the spacetime boson and spacetime fermion subspaces, as indicated in Table 3.1. Within these sub-blocks, there are further block diagonalizations arising from special properties of the one-loop form of the matrix elements of the perturbing Hamiltonian. For example, Fock space states built out of two bosonic creation operators that transform only under the internal $SO(4)$ mix only with themselves, thus providing a 16×16 dimensional diagonal sub-block. Within such sub-blocks, symmetry considerations are often sufficient to completely diagonalize the matrix or at least to reduce it to a low-dimensional diagonalization problem. In short, the problem reduces almost entirely to that of projecting the matrix elements of H_{int} on subspaces of the two-impurity Fock space defined by various symmetry properties. Determining the $SO(4) \times SO(4)$ symmetry labels of each eigenstate in the diagonalization will furthermore enable us to precisely match string states with gauge theory operators. In this subsection, we record for future reference the detailed arguments for the various special cases that must be dealt with in order to fully diagonalize the one-loop perturbation and characterize the irrep decomposition. Although the projections onto the various invariant subspaces are matters of simple algebra, that algebra is too complicated to be done by hand and we have resorted to symbolic manipulation programs. The end result of the diagonalization is quite simple and the reader willing to accept our results on faith can skip ahead to the end of this subsection.

We begin with a discussion of the action of the purely bosonic perturbation H_{BB} on the 64-dimensional Fock space created by pairs of bosonic creation operators. Part of this subspace connects via H_{BF} to the Fock space of spacetime bosons created by pairs of fermionic creation operators, and we will deal with it later. There is, however, a subspace that only connects to itself, through the purely bosonic perturbation H_{BB} . We will first deal with this purely bosonic block diagonalization, leading to

eigenvalues we will denote by Λ_{BB} . The eight bosonic modes lie in the $SO(4) \times SO(4)$ representations $(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$ and $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ (i.e., they are vectors in the $SO(4)$ subgroups descended from AdS_5 and S^5 , respectively). The key fact about H_{BB} is that the 16-dimensional spaces spanned by two $(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$ oscillators or by two $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ oscillators are closed under its action (it is also true that H_{BF} annihilates both of these subspaces). The $SO(4)$ representation content of the states created by such oscillator pairs is given by the formula $(\mathbf{2}, \mathbf{2}) \times (\mathbf{2}, \mathbf{2}) = (\mathbf{3}, \mathbf{3}) + (\mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}) + (\mathbf{1}, \mathbf{1})$ (we use $SU(2) \times SU(2)$ notation, rather than $SO(4)$, since it is unavoidable when we discuss fermions). By projecting the $O(\lambda')$ part of (3.5.6) onto these subspaces, one can directly read off the eigenvalues Λ_{BB} , with the results shown in Table 3.2. The identification of the representations associated with particular eigenvalues is easy to do on the basis of multiplicity. In any event, projection onto invariant subspaces is a simple matter of symmetrization or antisymmetrization of oscillator indices and can be done directly. The most important point to note is that the eigenvalues are successive even integers, a simple result and one that is consistent with our expectations from extended supersymmetry. It will be straightforward to match these states to gauge theory operators and compare eigenvalues with anomalous dimensions.

$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{BB}	$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{BB}
$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-6	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	2
$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$	-2	$(\mathbf{3}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-2
$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	-4	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	0

Table 3.2: Energy shifts at $O(1/J)$ for unmixed bosonic modes

The Fock space of spacetime bosons created by pairs of fermionic creation operators contains a similar pair of 16×16 diagonal sub-blocks. The construction and application of the relevant projection operators and the subsequent match-up with gauge theory operators is more complicated than on the bosonic side and we must develop some technical tools before we can obtain concrete results.

Just as H_{BB} is closed in the two 16-dimensional spaces of bosonic $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ or

$(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$ states, H_{FF} is closed on subspaces of bi-fermions spanned by a pair of $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ or a pair of $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ fermionic oscillators (i.e., $--$ or $++$ states, to use an obvious shorthand). The complete spectrum of eigenvalues from these subsectors of the Hamiltonian can be computed by projecting out the $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ and $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ spinors in H_{FF} (3.5.7). To do this, it will be helpful to express the eight-component spinors of the string theory in a basis which allows us to define fermionic oscillators labeled by their $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ and $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ representation content.

The original 32-component Majorana-Weyl spinors θ^I were reduced by the Weyl projection and a lightcone gauge condition to an eight-component spinor ψ^α (transforming in the 8_s of $SO(8)$). The generators of the four $SU(2)$ factors (3.5.9) of the manifest $SO(4) \times SO(4)$ symmetry can be expressed as 8×8 $SO(8)$ matrices as follows:

$$\begin{aligned} \Sigma_1^\pm &= -\frac{1}{4i}(\gamma^2\gamma^3 \pm \gamma^1\gamma^4) , & \Omega_1^\pm &= \frac{1}{4i}(-\gamma^6\gamma^7 \pm \gamma^5) , \\ \Sigma_2^\pm &= -\frac{1}{4i}(\gamma^3\gamma^1 \pm \gamma^2\gamma^4) , & \Omega_2^\pm &= \frac{1}{4i}(-\gamma^7\gamma^5 \pm \gamma^6) , \\ \Sigma_3^\pm &= -\frac{1}{4i}(\gamma^1\gamma^2 \pm \gamma^3\gamma^4) , & \Omega_3^\pm &= \frac{1}{4i}(-\gamma^5\gamma^6 \pm \gamma^7) . \end{aligned} \quad (3.5.17)$$

We will use the representation for the γ^A given in Appendix A (A.14) when we need to make these generators explicit. The 8_s spinor may be further divided into its $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ and $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ components $\hat{\psi}$ and $\tilde{\psi}$, respectively, and this suggests a useful basis change for the string creation operators: for the $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ spinor, we define four new objects w, x, y, z by

$$\hat{b}^\dagger = w \begin{pmatrix} 1 \\ 0 \\ 0 \\ -1 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix} + x \begin{pmatrix} 0 \\ 1 \\ 1 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix} + y \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \\ 1 \\ 0 \\ 0 \\ 0 \\ 1 \end{pmatrix} + z \begin{pmatrix} 0 \\ 1 \\ 1 \\ 0 \\ 0 \\ 1 \\ -1 \\ 0 \\ 0 \end{pmatrix} , \quad (3.5.18)$$

which we then organize in two different ways into two-component complex spinors:

$$\begin{aligned}\zeta &= \begin{pmatrix} w + iy \\ z + ix \end{pmatrix}, & \varphi &= \begin{pmatrix} -z + ix \\ w - iy \end{pmatrix} \Leftarrow \Sigma_i^-, \\ \bar{\zeta} &= \begin{pmatrix} w + iy \\ -z + ix \end{pmatrix}, & \bar{\varphi} &= \begin{pmatrix} z + ix \\ w - iy \end{pmatrix} \Leftarrow \Omega_i^-. \end{aligned} \quad (3.5.19)$$

This organization into two-spinors is meant to show how components of $\hat{\psi}$ transform under the two $SU(2)$ factors that act nontrivially on them. As may be verified from the explicit forms of the $SU(2)$ generators obtained by substituting (A.14) into (3.5.17), the two-component spinors ζ and φ transform as $(\mathbf{1}, \mathbf{2})$ under the first $SO(4)$ and the spinors $\bar{\zeta}$ and $\bar{\varphi}$ transform as $(\mathbf{1}, \mathbf{2})$ under the second $SO(4)$ of $SO(4) \times SO(4)$. The explicit realization of the two $SU(2)$ factors involved here is found in this way to be

$$\begin{aligned}\Sigma_1^- &= \begin{pmatrix} 0 & 1/2 \\ 1/2 & 0 \end{pmatrix}, & \Omega_1^- &= \begin{pmatrix} 0 & 1/2 \\ 1/2 & 0 \end{pmatrix}, \\ \Sigma_2^- &= \begin{pmatrix} 0 & i/2 \\ -i/2 & 0 \end{pmatrix}, & \Omega_2^- &= \begin{pmatrix} 0 & -i/2 \\ i/2 & 0 \end{pmatrix}, \\ \Sigma_3^- &= \begin{pmatrix} 1/2 & 0 \\ 0 & -1/2 \end{pmatrix}, & \Omega_3^- &= \begin{pmatrix} 1/2 & 0 \\ 0 & -1/2 \end{pmatrix}. \end{aligned} \quad (3.5.20)$$

One may similarly decompose $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ spinors and express the corresponding

generators Σ^+ and Ω^+ . We decompose $\tilde{\psi}$ into components $\bar{w}, \bar{x}, \bar{y}, \bar{z}$ according to

$$\tilde{b}^\dagger = \bar{w} \begin{pmatrix} 1 \\ 0 \\ 0 \\ 1 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix} + \bar{x} \begin{pmatrix} 0 \\ 1 \\ -1 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix} + \bar{y} \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \\ 1 \\ 0 \\ 0 \\ -1 \end{pmatrix} + \bar{z} \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 1 \\ 1 \\ 0 \end{pmatrix}, \quad (3.5.21)$$

and rearrange them into two-component complex spinors:

$$\begin{aligned} \xi &= \begin{pmatrix} \bar{z} + i\bar{x} \\ \bar{w} + i\bar{y} \end{pmatrix}, & \eta &= \begin{pmatrix} \bar{w} - i\bar{y} \\ -\bar{z} + i\bar{x} \end{pmatrix} \Leftarrow \Sigma_i^+, \\ \bar{\xi} &= \begin{pmatrix} -\bar{z} + i\bar{x} \\ \bar{w} + i\bar{y} \end{pmatrix}, & \bar{\eta} &= \begin{pmatrix} \bar{w} - i\bar{y} \\ \bar{z} + i\bar{x} \end{pmatrix} \Leftarrow \Omega_i^+. \end{aligned} \quad (3.5.22)$$

The corresponding explicit $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ generators are given by

$$\begin{aligned} \Sigma_1^+ &= \begin{pmatrix} 0 & -1/2 \\ -1/2 & 0 \end{pmatrix}, & \Omega_1^+ &= \begin{pmatrix} 0 & 1/2 \\ 1/2 & 0 \end{pmatrix}, \\ \Sigma_2^+ &= \begin{pmatrix} 0 & i/2 \\ -i/2 & 0 \end{pmatrix}, & \Omega_2^+ &= \begin{pmatrix} 0 & -i/2 \\ i/2 & 0 \end{pmatrix}, \\ \Sigma_3^+ &= \begin{pmatrix} 1/2 & 0 \\ 0 & -1/2 \end{pmatrix}, & \Omega_3^+ &= \begin{pmatrix} 1/2 & 0 \\ 0 & -1/2 \end{pmatrix}. \end{aligned} \quad (3.5.23)$$

These observations will make it possible to construct linear combinations of products of components of ψ^α transforming in chosen irreps of $SO(4) \times SO(4)$.

Let us now use this machinery to analyze the perturbation matrix on spacetime

bosons created by two fermionic creation operators (bi-fermions). As explained in the discussion of (3.5.7), H_{FF} is block-diagonal on the 16-dimensional $++$ or $--$ bi-fermionic subspaces. To project out the $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $++$ block of H_{FF} , we simply act on all indices of (3.5.7) with the Π_+ projection operator:

$$\begin{aligned} \left\langle J \left| \tilde{b}_n^\alpha \tilde{b}_{-n}^\beta (H_{\text{FF}}) \tilde{b}_{-n}^{\gamma\dagger} \tilde{b}_n^{\delta\dagger} \right| J \right\rangle &= -2 \frac{n^2 \lambda'}{J} \Pi_+^{\alpha\delta} \Pi_+^{\beta\gamma} + \frac{n^2 \lambda'}{24J} \left\{ \left[(\Pi_+ \gamma^{ij} \Pi_+)^{\alpha\delta} (\Pi_+ \gamma^{ij} \Pi_+)^{\beta\gamma} \right. \right. \\ &\quad \left. \left. + (\Pi_+ \gamma^{ij} \Pi_+)^{\alpha\beta} (\Pi_+ \gamma^{ij} \Pi_+)^{\gamma\delta} - (\Pi_+ \gamma^{ij} \Pi_+)^{\alpha\gamma} (\Pi_+ \gamma^{ij} \Pi_+)^{\beta\delta} \right] \right. \\ &\quad \left. - \left[(\Pi_+ \gamma^{i'j'} \Pi_+)^{\alpha\delta} (\Pi_+ \gamma^{i'j'} \Pi_+)^{\beta\gamma} + (\Pi_+ \gamma^{i'j'} \Pi_+)^{\alpha\beta} (\Pi_+ \gamma^{i'j'} \Pi_+)^{\gamma\delta} \right. \right. \\ &\quad \left. \left. - (\Pi_+ \gamma^{i'j'} \Pi_+)^{\alpha\gamma} (\Pi_+ \gamma^{i'j'} \Pi_+)^{\beta\delta} \right] \right\}. \end{aligned} \quad (3.5.24)$$

The $SO(4) \times SO(4)$ representation content of this subspace is specified by $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) \times (\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) = (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1}) \oplus (\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) \oplus (\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) \oplus (\mathbf{3}, \mathbf{1}; \mathbf{3}, \mathbf{1})$ and we must further project onto individual irreducible representations in order to identify the eigenvalues.

With the tools we have built up in the last few paragraphs, we are in a position to directly project out some of the desired irreducible representations. Bi-fermions of $++$ type transforming as scalars under the first $SO(4)$ (i.e., under Σ_i^+) are constructed by making $SU(2)$ invariants out of the two-component spinors ξ and η . There are four such objects:

$$\begin{aligned} \xi_{-n} \tau_2 \xi_n, & \quad \xi_{-n} \tau_2 \eta_n, \\ \eta_{-n} \tau_2 \xi_n, & \quad \eta_{-n} \tau_2 \eta_n, \end{aligned} \quad (3.5.25)$$

where τ_2 is the second Pauli matrix. At the same time, they must also comprise a $\mathbf{3}$ and a $\mathbf{1}$ under the second $SO(4)$ (i.e., under Ω_i^+). To identify the irreducible linear combinations, one has to re-express the objects in (3.5.25) in terms of the spinors $\bar{\xi}$ and $\bar{\eta}$ that transform simply under Ω_i^+ . Representative results for properly normalized

creation operators of $++$ bi-fermion states in particular $SO(4) \times SO(4)$ irreps are

$$\begin{aligned}
& -\frac{1}{2} (\xi_{-n} \tau_2 \eta_n - \eta_{-n} \tau_2 \xi_n) \quad (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1}) , \quad \Lambda_{\text{FF}} = -2 , \\
& \left. \begin{aligned} & \frac{1}{2} (\xi_{-n} \tau_2 \eta_n + \eta_{-n} \tau_2 \xi_n) \\ & \frac{i}{2} (\xi_{-n} \tau_2 \xi_n + \eta_{-n} \tau_2 \eta_n) \\ & -\frac{1}{2} (\xi_{-n} \tau_2 \xi_n - \eta_{-n} \tau_2 \eta_n) \end{aligned} \right\} (\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) , \quad \Lambda_{\text{FF}} = 0 . \quad (3.5.26)
\end{aligned}$$

We simply have to re-express the ξ, η bilinears in terms of the original spinor creation operators \tilde{b} in order to obtain an explicit projection of the matrix elements (3.5.24) onto irreducible subspaces and to obtain the eigenvalues Λ_{FF} associated with each irrep. A parallel analysis of states constructed by forming normalized $SU(2)$ invariants from $\bar{\xi}$ and $\bar{\eta}$ gives another irrep and eigenvalue:

$$\begin{aligned}
& \left. \begin{aligned} & \frac{1}{2} (\bar{\xi}_{-n} \tau_2 \bar{\eta}_n + \bar{\eta}_{-n} \tau_2 \bar{\xi}_n) \\ & \frac{i}{2} (\bar{\xi}_{-n} \tau_2 \bar{\xi}_n + \bar{\eta}_{-n} \tau_2 \bar{\eta}_n) \\ & -\frac{1}{2} (\bar{\xi}_{-n} \tau_2 \bar{\xi}_n - \bar{\eta}_{-n} \tau_2 \bar{\eta}_n) \end{aligned} \right\} (\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) , \quad \Lambda_{\text{FF}} = -4 . \quad (3.5.27)
\end{aligned}$$

By similar arguments, whose details we will omit, one can construct the creation operator for the normalized $(\mathbf{3}, \mathbf{1}; \mathbf{3}, \mathbf{1})$ or $++$ bi-fermion and find the eigenvalue $\Lambda_{\text{FF}} = -2$.

An exactly parallel analysis of $\left\langle J \left| \hat{b} \hat{b} (H_{\text{FF}}) \hat{b}^\dagger \hat{b}^\dagger \right| J \right\rangle$ on the 16-dimensional subspace spanned by $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ bi-fermions yields the same eigenvalue spectrum. The creation operators of irreducible states (built this time out of ζ and ϕ) and their eigenvalues are

$$-\frac{1}{2} (\zeta_{-n} \tau_2 \varphi_n - \varphi_{-n} \tau_2 \zeta_n) \quad (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1}) , \quad \Lambda_{\text{FF}} = -2 ,$$

$$\left. \begin{aligned} & \frac{1}{2} (\zeta_{-n} \tau_2 \varphi_n + \varphi_{-n} \tau_2 \zeta_n) \\ & \frac{i}{2} (\zeta_{-n} \tau_2 \zeta_n + \varphi_{-n} \tau_2 \varphi_n) \\ & -\frac{1}{2} (\zeta_{-n} \tau_2 \zeta_n - \varphi_{-n} \tau_2 \varphi_n) \end{aligned} \right\} (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3}) , \quad \Lambda_{\text{FF}} = 0 , \quad (3.5.28)$$

$$\left. \begin{aligned} & \frac{1}{2} (\bar{\zeta}_{-n} \tau_2 \bar{\varphi}_n + \bar{\varphi}_{-n} \tau_2 \bar{\zeta}_n) \\ & \frac{i}{2} (\bar{\zeta}_{-n} \tau_2 \bar{\zeta}_n + \bar{\varphi}_{-n} \tau_2 \bar{\varphi}_n) \\ & -\frac{1}{2} (\bar{\zeta}_{-n} \tau_2 \bar{\zeta}_n - \bar{\varphi}_{-n} \tau_2 \bar{\varphi}_n) \end{aligned} \right\} (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1}) , \quad \Lambda_{\text{FF}} = -4 . \quad (3.5.29)$$

The overall results for this sector are displayed in Table 3.3.

$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{FF}	$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{FF}
$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-2	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-2
$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1})$	0	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	0
$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-4	$(\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-4
$(\mathbf{3}, \mathbf{1}; \mathbf{3}, \mathbf{1})$	-2	$(\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{3})$	-2

Table 3.3: Energy shifts of states created by two fermions in $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$

To this point, we have been able to study specific projections of the H_{BB} and H_{FF} subsectors by choosing states that are not mixed by H_{BF} . We now must deal with the subspace of spacetime boson two-impurity states that is not annihilated by H_{BF} . This 64-dimensional space is spanned by pairs of bosonic creation operators taken from different $SO(4)$ subgroups and pairs of fermionic creation operators of opposite Π -parity. The representation content of these creation-operator pairs is such that the states in this sector all belong to $(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$ irreps. This space is of course also acted on by H_{BB} and H_{FF} , so we will need the matrix elements of all three pieces of the Hamiltonian as they act on this subspace. By applying the appropriate projections to the general one-loop matrix elements, we obtain the expressions

$$\left\langle J \left| a_n^A a_{-n}^B (H_{\text{BB}}) a_{-n}^{C\dagger} a_n^{D\dagger} \right| J \right\rangle \rightarrow -2 \frac{n^2 \lambda'}{J} \left(\delta^{ad'} \delta^{b'c} + \delta^{a'd} \delta^{bc'} + \delta^{ad} \delta^{b'c'} + \delta^{a'd'} \delta^{bc} \right) , \quad (3.5.30)$$

$$\begin{aligned}
\left\langle J \left| b_n^{\alpha} b_{-n}^{\beta} (H_{\text{BF}}) a_{-n}^{A\dagger} a_n^{B\dagger} \right| J \right\rangle &\rightarrow \frac{n^2 \lambda'}{2J} \left[\left(\Pi_+ \gamma^{ab'} \Pi_- \right)^{\alpha\beta} - \left(\Pi_+ \gamma^{a'b} \Pi_- \right)^{\alpha\beta} \right. \\
&\quad \left. + \left(\Pi_- \gamma^{ab'} \Pi_+ \right)^{\alpha\beta} - \left(\Pi_- \gamma^{a'b} \Pi_+ \right)^{\alpha\beta} \right], \tag{3.5.31}
\end{aligned}$$

$$\begin{aligned}
\left\langle J \left| b_n^{\alpha} b_{-n}^{\beta} (H_{\text{FF}}) b_{-n}^{\gamma\dagger} b_n^{\delta\dagger} \right| J \right\rangle &\rightarrow -2 \frac{n^2 \lambda'}{J} \left(\Pi_+^{\alpha\delta} \Pi_-^{\beta\gamma} + \Pi_-^{\alpha\delta} \Pi_+^{\beta\gamma} \right) \\
&+ \frac{n^2 \lambda'}{24J} \left\{ \left[\left(\Pi_+ \gamma^{ij} \Pi_+ \right)^{\alpha\delta} \left(\Pi_- \gamma^{ij} \Pi_- \right)^{\beta\gamma} + \left(\Pi_+ \gamma^{ij} \Pi_- \right)^{\alpha\beta} \left(\Pi_- \gamma^{ij} \Pi_+ \right)^{\gamma\delta} \right. \right. \\
&\quad \left. \left. - \left(\Pi_+ \gamma^{ij} \Pi_- \right)^{\alpha\gamma} \left(\Pi_- \gamma^{ij} \Pi_+ \right)^{\beta\delta} \right] - \left[\left(\Pi_+ \gamma^{i'j'} \Pi_+ \right)^{\alpha\delta} \left(\Pi_- \gamma^{i'j'} \Pi_- \right)^{\beta\gamma} \right. \right. \\
&\quad \left. \left. + \left(\Pi_+ \gamma^{i'j'} \Pi_- \right)^{\alpha\beta} \left(\Pi_- \gamma^{i'j'} \Pi_+ \right)^{\gamma\delta} - \left(\Pi_+ \gamma^{i'j'} \Pi_- \right)^{\alpha\gamma} \left(\Pi_- \gamma^{i'j'} \Pi_+ \right)^{\beta\delta} \right] \right. \\
&\quad \left. + \left[\left(\Pi_- \gamma^{ij} \Pi_- \right)^{\alpha\delta} \left(\Pi_+ \gamma^{ij} \Pi_+ \right)^{\beta\gamma} + \left(\Pi_- \gamma^{ij} \Pi_+ \right)^{\alpha\beta} \left(\Pi_+ \gamma^{ij} \Pi_- \right)^{\gamma\delta} \right. \right. \\
&\quad \left. \left. - \left(\Pi_- \gamma^{ij} \Pi_+ \right)^{\alpha\gamma} \left(\Pi_+ \gamma^{ij} \Pi_- \right)^{\beta\delta} \right] - \left[\left(\Pi_- \gamma^{i'j'} \Pi_- \right)^{\alpha\delta} \left(\Pi_+ \gamma^{i'j'} \Pi_+ \right)^{\beta\gamma} \right. \right. \\
&\quad \left. \left. + \left(\Pi_- \gamma^{i'j'} \Pi_+ \right)^{\alpha\beta} \left(\Pi_+ \gamma^{i'j'} \Pi_- \right)^{\gamma\delta} - \left(\Pi_- \gamma^{i'j'} \Pi_+ \right)^{\alpha\gamma} \left(\Pi_+ \gamma^{i'j'} \Pi_- \right)^{\beta\delta} \right] \right\}. \tag{3.5.32}
\end{aligned}$$

Since the 64-dimensional space must contain four copies of the $(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$ irrep, the diagonalization problem is really only 4×4 and quite easy to solve. The results for the eigenvalues appear in Table 3.4. Collecting the above results, we present the

$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{BF}
$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	-4
$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2}) \times 2$	-2
$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	0

Table 3.4: String eigenstates in the subspace for which H_{BF} has non-zero matrix elements

complete $SO(4)_{AdS} \times SO(4)_{S^5}$ decomposition of spacetime boson two-impurity states in Table 3.5.

By projecting out closed subspaces of the one-loop Hamiltonian we have success-

	$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ
H_{BB}	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-6
	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	2
	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	-4
	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	0
	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$	-2
	$(\mathbf{3}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-2
	$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ
H_{FF}	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-2
	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-2
	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	0
	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-4
	$(\mathbf{3}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{3})$	-2
H_{BF}	$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	0
	$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2}) \times 2$	-2
	$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	-4

Table 3.5: Group decomposition of the 128 two-impurity spacetime bosons

fully classified each of the energy levels in the bosonic Fock space with an $SO(4) \times SO(4)$ symmetry label. Similar arguments can be applied to the fermionic Fock space, where two-impurity string states mix individual bosonic and fermionic oscillators (we omit the details). A summary of these results for all states, including spacetime fermions, is given in Table 3.6. The important fact to note is that the Λ eigenvalues and their multiplicities are exactly as required for consistency with the full $PSU(2, 2|4)$ symmetry of the theory. This is a nontrivial result since the quantization procedure does not make the full symmetry manifest. It is also a very satisfying check of the overall correctness of the extremely complicated set of procedures we were forced to use. We can now proceed to a comparison with gauge theory anomalous dimensions.

Level	0	2	4	6	8
Mult.	1	28	70	28	1
Λ_{Bose}	-6	-4	-2	0	2

Level	1	3	5	7
Mult.	8	56	56	8
Λ_{Fermi}	-5	-3	-1	1

Table 3.6: First-order energy shift summary: complete two-impurity string multiplet

3.5.6 Gauge theory comparisons

The most comprehensive analysis of one-loop anomalous dimensions of BMN operators and their organization into supersymmetry multiplets was given in [28]. As stated in [26], the above string theory calculations are in perfect agreement with the one-loop gauge theory predictions. For completeness, we present a summary of the spectrum of dimensions of gauge theory operators along with a sampling of information about their group transformation properties.

The one-loop formula for operator dimensions takes the generic form

$$\Delta_n^R = 2 + \frac{g_{YM}^2 N_c}{R^2} n^2 \left(1 + \frac{\bar{\Lambda}}{R} + O(R^{-2}) \right) . \quad (3.5.33)$$

The $O(R^{-1})$ correction $\bar{\Lambda}$ for the set of two-impurity operators is predicted to match the corresponding $O(J^{-1})$ energy correction to two-impurity string states, labeled above by Λ . Part of the motivation for performing the special projections on two-impurity string states detailed above was to emerge with specific symmetry labels for each of the string eigenstates. String states of a certain representation content of the residual $SO(4) \times SO(4)$ symmetry of $AdS_5 \times S^5$ are expected, by duality, to map to gauge theory operators with the same representation labels in the $SL(2, \mathbf{C})$ Lorentz and $SU(4)$ R -charge sectors of the gauge theory. Knowing the symmetry content of the string eigenstates therefore allows us to test this mapping in detail.

The bosonic sector of the gauge theory, characterized by single-trace operators with two bosonic insertions in the trace, appears in Table 3.7. The set of operators comprising Lorentz scalars clearly agrees with the corresponding pure-boson string states in Table 3.5, which are scalars in AdS_5 . Operators containing pairs of spacetime derivatives correspond to string theory states that are scalars of the S^5 subspace. The bi-fermion sector of the string theory corresponds to the set of two-gluino operators in the gauge theory. A few of these operators are listed in Table 3.8. These states, which form either spacetime or R -charge scalars, clearly agree with their string theory counterparts, which were constructed explicitly above. The string states appearing in

Operator	$SO(4)_{AdS} \times SO(4)_{S^5}$	$\bar{\Lambda}$
$\Sigma_A \text{tr} (\phi^A Z^p \phi^A Z^{R-p})$	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-6
$\text{tr} (\phi^{(i} Z^p \phi^{j)} Z^{R-p})$	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$	-2
$\text{tr} (\phi^{[i} Z^p \phi^{j]} Z^{R-p})$	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	-4
$\text{tr} (\nabla_\mu Z Z^p \nabla^\mu Z Z^{R-2-p})$	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	2
$\text{tr} (\nabla_{(\mu} Z Z^p \nabla_{\nu)} Z Z^{R-2-p})$	$(\mathbf{3}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-2
$\text{tr} (\nabla_{[\mu} Z Z^p \nabla_{\nu]} Z Z^{R-2-p})$	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	0

Table 3.7: Bosonic gauge theory operators: either spacetime or R -charge singlet.

the $(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$ representation (listed in Table 3.4) correspond to the operators listed in Table 3.9. Finally, the complete supermultiplet spectrum of two-impurity gauge theory operators appears in Table 3.10. The extended supermultiplet spectrum is in perfect agreement with the complete one-loop string theory spectrum in Table 3.6 above.

Operator	$SO(4)_{AdS} \times SO(4)_{S^5}$	$\bar{\Lambda}$
$\text{tr} (\chi^{[\alpha} Z^p \chi^{\beta]} Z^{R-1-p})$	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-2
$\text{tr} (\chi^{(\alpha} Z^p \chi^{\beta)} Z^{R-1-p})$	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1})$	0
$\text{tr} (\chi [\sigma_\mu, \tilde{\sigma}_\nu] Z^p \chi Z^{R-1-p})$	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-4

Table 3.8: Bosonic gauge theory operators with two gluino impurities.

Operator	$SO(4)_{AdS} \times SO(4)_{S^5}$	$\bar{\Lambda}$
$\text{tr} (\phi^i Z^p \nabla_\mu Z Z^{R-1-p}) + \dots$	$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	-4
$\text{tr} (\phi^i Z^p \nabla_\mu Z Z^{R-1-p})$	$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	-2
$\text{tr} (\phi^i Z^p \nabla_\mu Z Z^{R-1-p}) + \dots$	$(\mathbf{2}, \mathbf{2}; \mathbf{2}, \mathbf{2})$	0

Table 3.9: Bosonic gauge theory operators: spacetime and R -charge non-singlets

Level	0	1	2	3	4	5	6	7	8
Multiplicity	1	8	28	56	70	56	28	8	1
$\delta E \times (R^2/g_{YM}^2 N_c n^2)$	$-6/R$	$-5/R$	$-4/R$	$-3/R$	$-2/R$	$-1/R$	0	$1/R$	$2/R$

Table 3.10: Anomalous dimensions of two-impurity operators

3.6 Energy spectrum at all loops in λ'

To make comparisons with gauge theory dimensions at one loop in $\lambda = g_{YM}^2 N_c$, we have expanded all string energies in powers of the modified 't Hooft coupling $\lambda' = g_{YM}^2 N_c / R^2$. The string theory analysis is exact to all orders in λ' , however, and it is possible to extract a formula for the $O(1/J)$ string energy corrections that is exact in λ' and suitable for comparison with higher-order corrections to operator dimensions in the gauge theory. In practice, it is slightly more difficult to diagonalize the string Hamiltonian when the matrix elements are not expanded in small λ' . This is mainly because, beyond leading order, H_{BF} acquires additional terms that mix bosonic indices in the same $SO(4)$ and also mix bi-fermionic indices in the same $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ or $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ representation. Instead of a direct diagonalization of the entire 128-dimensional subspace of spacetime bosons, for example, we find it more convenient to exploit the ‘dimension reduction’ that can be achieved by projecting the full Hamiltonian onto individual irreps.

For example, the $(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$ irrep appears four times in Table 3.5 and is present at levels $L = 0, 4, 8$ in the supermultiplet. To get the exact eigenvalues for this irrep, we will have to diagonalize a 4×4 matrix. The basis vectors of this bosonic sector comprise singlets of the two $SO(4)$ subgroups $(a^{\dagger a} a^{\dagger a} |J\rangle)$ and $(a^{\dagger a'} a^{\dagger a'} |J\rangle)$ plus two bi-fermion singlets constructed from the $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ and $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ creation operators $(\hat{b}^{\dagger \alpha} \hat{b}^{\dagger \alpha} |J\rangle)$ and $(\tilde{b}^{\dagger \alpha} \tilde{b}^{\dagger \alpha} |J\rangle)$. The different Hamiltonian matrix elements that enter the 4×4 matrix are symbolically indicated in Table 3.11. It is a simple matter to

project the general expressions for matrix elements of H_{BB} , etc., onto singlet states and so obtain the matrix as an explicit function of λ', n . The matrix can be exactly diagonalized and yields the following energies:

$$\begin{aligned}
E_0(n, J) &= 2\sqrt{1 + \lambda' n^2} - \frac{n^2 \lambda'}{J} \left[2 + \frac{4}{\sqrt{1 + n^2 \lambda'}} \right] + O(1/J^2) , \\
E_4(n, J) &= 2\sqrt{1 + \lambda' n^2} - \frac{2n^2 \lambda'}{J} + O(1/J^2) , \\
E_8(n, J) &= 2\sqrt{1 + \lambda' n^2} - \frac{n^2 \lambda'}{J} \left[2 - \frac{4}{\sqrt{1 + n^2 \lambda'}} \right] + O(1/J^2) . \quad (3.6.1)
\end{aligned}$$

The subscript $L = 0, 4, 8$ indicates the supermultiplet level to which the eigenvalue connects in the weak coupling limit. The middle eigenvalue ($L = 4$) is doubly degenerate, as it was in the one-loop limit.

H_{int}	$a^{\dagger a} a^{\dagger a} J\rangle$	$a^{\dagger a'} a^{\dagger a'} J\rangle$	$\hat{b}^{\dagger \alpha} \hat{b}^{\dagger \alpha} J\rangle$	$\tilde{b}^{\dagger \alpha} \tilde{b}^{\dagger \alpha} J\rangle$
$\langle J a^a a^a$	H_{BB}	H_{BB}	H_{BF}	H_{BF}
$\langle J a^{a'} a^{a'}$	H_{BB}	H_{BB}	H_{BF}	H_{BF}
$\langle J \hat{b}^\alpha \hat{b}^\alpha$	H_{BF}	H_{BF}	H_{FF}	H_{FF}
$\langle J \tilde{b}^\alpha \tilde{b}^\alpha$	H_{BF}	H_{BF}	H_{FF}	H_{FF}

Table 3.11: Singlet projection at finite λ'

There are two independent 2×2 matrices that mix states at levels $L = 2, 6$. According to Table 3.5, one can project out the antisymmetric bosonic and antisymmetric bi-fermionic states in the irrep $(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$ or in the irrep $(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$. The results of eqns. (3.5.26, 3.5.27, 3.5.28, 3.5.29) can be used to carry out the needed projections and obtain explicit forms for the matrix elements of the perturbing Hamiltonian. The actual 2×2 diagonalization is trivial to do and both problems give the same result. The final result for the energy levels

(using the same notation as before) is

$$\begin{aligned} E_2(n, J) &= 2\sqrt{1 + \lambda'n^2} - \frac{n^2\lambda'}{J} \left[2 + \frac{2}{\sqrt{1 + n^2\lambda'}} \right] + O(1/J^2) , \\ E_6(n, J) &= 2\sqrt{1 + \lambda'n^2} - \frac{n^2\lambda'}{J} \left[2 - \frac{2}{\sqrt{1 + n^2\lambda'}} \right] + O(1/J^2) . \end{aligned} \quad (3.6.2)$$

We can carry out similar diagonalizations for the remaining irreps of Table 3.5, but no new eigenvalues are encountered: the energies already listed are the exact energies of the $L = 0, 2, 4, 6, 8$ levels. It is also easy to see that the degeneracy structure of the exact levels is the same as the one-loop degeneracy.

The odd levels of the supermultiplet are populated by the 128-dimensional space-time fermions, and this sector of the theory can be diagonalized directly. Proceeding in a similar fashion as in the bosonic sector, we find exact energy eigenvalues for the $L = 1, 3, 5, 7$ levels (with unchanged multiplicities). We refrain from stating the individual results because the entire supermultiplet spectrum, bosonic and fermionic, can be written in terms of a single concise formula: to leading order in $1/J$ and all orders in λ' , the energies of the two-impurity multiplet are given by

$$E_L(n, J) = 2\sqrt{1 + \lambda'n^2} - \frac{n^2\lambda'}{J} \left[2 + \frac{(4-L)}{\sqrt{1 + n^2\lambda'}} \right] + O(1/J^2) , \quad (3.6.3)$$

where $L = 0, 1, \dots, 8$ indicates the level within the supermultiplet. The degeneracies and irrep content are identical to what we found at one loop in λ' . This expression can be rewritten, correct to order J^{-2} , as follows:

$$E_L(n, J) \approx 2\sqrt{1 + \frac{\lambda n^2}{(J - L/2)^2}} - \frac{n^2\lambda}{(J - L/2)^3} \left[2 + \frac{4}{\sqrt{1 + \lambda n^2/(J - L/2)^2}} \right] . \quad (3.6.4)$$

This shows that, within this expansion, the joint dependence on J and L is exactly what is required for extended supersymmetry multiplets. This is a rather nontrivial functional requirement, and a stringent check on the correctness of our quantization

procedure (independent of any comparison with gauge theory).

In order to make contact with gauge theory we expand (3.6.3) in λ' , obtaining

$$E_L(n, J) \approx \left[2 + \lambda' n^2 - \frac{1}{4}(\lambda' n^2)^2 + \frac{1}{8}(\lambda' n^2)^3 + \dots \right] + \frac{1}{J} \left[n^2 \lambda' (L - 6) + (n^2 \lambda')^2 \left(\frac{4 - L}{2} \right) + (n^2 \lambda')^3 \left(\frac{3L - 12}{8} \right) + \dots \right]. \quad (3.6.5)$$

We can now address the comparison with higher-loop results on gauge theory operator dimensions. Beisert, Kristjansen and Staudacher [31] computed the two-loop correction to the anomalous dimensions of a convenient class of operators lying at level four in the supermultiplet. The operators in question lie in a symmetric-traceless irrep of an $SO(4)$ subgroup of the R -charge and are guaranteed by group theory not to mix with any other fields [31]. The following expression for the two-loop anomalous dimension was found:

$$\delta\Delta_n^R = -\frac{g_{YM}^4 N_c^2}{\pi^4} \sin^4 \frac{n\pi}{R+1} \left(\frac{1}{4} + \frac{\cos^2 \frac{n\pi}{R+1}}{R+1} \right). \quad (3.6.6)$$

As explained above, $\mathcal{N} = 4$ supersymmetry ensures that the dimensions of operators at other levels of the supermultiplet will be obtained by making the substitution $R \rightarrow R + 2 - L/2$ in the expression for the dimension of the $L = 4$ operator. Making that substitution and taking the large- R limit we obtain a general formula for the two-loop, large- R correction to the anomalous dimension of the general two-impurity operator:

$$\begin{aligned} \delta\Delta_n^{R,L} &= -\frac{g_{YM}^4 N_c^2}{\pi^4} \sin^4 \frac{n\pi}{R+3-L/2} \left(\frac{1}{4} + \frac{\cos^2 \frac{n\pi}{R+3-L/2}}{R+3-L/2} \right) \\ &\approx -\frac{1}{4}(\lambda' n^2)^2 + \frac{1}{2}(\lambda' n^2)^2 \frac{4-L}{R} + O(1/R^2), \end{aligned} \quad (3.6.7)$$

Using the identification $R \rightleftharpoons J$ specified by duality, we see that this expression matches the corresponding string result in (3.6.5) to $O(1/J)$, confirming the AdS/CFT correspondence to two loops in the gauge coupling.

The three-loop correction to the dimension of this same class of $L = 4$ gauge theory operators has recently been definitively determined [33]. The calculation involves a remarkable interplay between gauge theory and integrable spin chain models [30, 31, 69, 89]. The final result is

$$\delta\Delta_n^R = \left(\frac{\lambda}{\pi^2}\right)^3 \sin^6 \frac{n\pi}{R+1} \left[\frac{1}{8} + \frac{\cos^2 \frac{n\pi}{R+1}}{4(R+1)^2} \left(3R + 2(R+6) \cos^2 \frac{n\pi}{R+1} \right) \right]. \quad (3.6.8)$$

If we apply to this expression the same logic applied to the two-loop gauge theory result (3.6.6), we obtain the following three-loop correction to the anomalous dimension of the general level of the two-impurity operator supermultiplet:

$$\delta\Delta_n^{R,L} \approx \frac{1}{8}(\lambda'n^2)^3 - \frac{1}{8}(\lambda'n^2)^2 \frac{8-3L}{R} + O(1/R^2). \quad (3.6.9)$$

We see that this expression differs from the third-order contribution to the string result (3.6.5) for the corresponding quantity. The difference is a constant shift and one might hope to absorb it in a normal-ordering constant. However, our discussion of the normal-ordering issue earlier in the chapter seems to exclude any such freedom.

3.7 Discussion

In this chapter we have given a detailed account of the quantization of the first curvature correction to type IIB superstring theory in the plane-wave limit of $AdS_5 \times S^5$. We have presented the detailed diagonalization of the resulting perturbing Hamiltonian on the degenerate subspace of two-impurity states, obtaining string energy corrections that can be compared with higher-loop anomalous dimensions of gauge theory operators. Beyond the Penrose limit, the holographic mapping between each side of the correspondence is intricate and nontrivial, and works perfectly to two loops in the gauge coupling. The agreement, however, appears to break down at three loops. (Similar three-loop disagreements have appeared in semiclassical string analyses; see,

for example, ref. [37].) This troubling issue was first observed in [26], at which time the third-order gauge theory anomalous dimension was somewhat conjectural. In the intervening time, the third-order result (3.6.8) acquired a solid basis, thus confirming the mismatch. Several questions arise about this mismatch: is it due to a failure of the AdS/CFT correspondence itself? Does it signal the need to modify the world-sheet string action? Is it simply that the perturbative approach to the gauge theory anomalous dimensions is not adequate in the relevant limits?

Regarding this final point, a specific explanation has been proposed that may account for the disagreement with gauge theory at three loops. The essential idea is that certain types of gauge theory mixing terms that connect fields within single-trace operators are dropped in the particular limit that is taken in this setup. This amounts to a plausible order-of-limits problem: we will leave a more detailed discussion of this proposal for Chapter 7. Despite vigorous investigation from several directions, all of these questions remain open.

Chapter 4

The curvature expansion: Three impurities

Thus far, we have seen that attempts to push the original results of BMN further have gone in two independent directions. In the gauge theory, the calculation of anomalous dimensions of BMN operators has been greatly simplified by Minahan and Zarembo's discovery that the problem can be mapped to that of computing the energies of certain integrable spin chains [39]. Based on this development, calculations in certain sectors of the theory have been carried out to three loops in the 't Hooft coupling λ [31,33].¹ Furthermore, we have shown in the previous chapter that the quantization of the GS string in the $AdS_5 \times S^5$ background has developed far enough to enable perturbative computations of the effect of worldsheet interactions on the spectrum of the string when it is boosted to large, but finite, angular momentum J [26,27,79]. As demonstrated in Chapter 3, these two approaches lead to different expansions of operator anomalous dimensions (or string eigenenergies): on the gauge theory side, one naturally has an expansion in the coupling constant λ that is typically exact in R -charge; on the string theory side one has an expansion in inverse powers of angular momentum J (the dual of gauge theory R -charge) that is exact in λ .

The expansion on the string side is difficult and has so far been carried out to $O(1/J)$ for two-impurity states (i.e., states with two string oscillators excited). The

¹We note that the conjectural three-loop computation of [31] was solidified by field theoretic methods in [33].

resulting functions of the loop expansion parameter λ can be compared with the large R -charge expansion of two-impurity BMN operators in the gauge theory to provide new and stringent tests of the AdS/CFT correspondence. As mentioned above, gauge theory technology has made it possible to compute anomalous dimensions of certain two-impurity BMN operators out to at least three-loop order. The agreement between dual quantities is perfect out to two-loop order but, surprisingly, seems to break down at three loops [26, 27]. Exactly what this means for the AdS/CFT correspondence is not yet clear but, given the circumstances, it seems appropriate to at least look for further data on the disagreement in the hope of finding some instructive systematics. The subject of this chapter is to pursue one possible line of attack in which we extend the calculations described above to higher-impurity string states and gauge theory operators. The extension of our two-impurity results to higher impurities is not a straightforward matter on either side of the correspondence and gets more complex as the number of impurities increases. We focus here on the three-impurity case, where we obtain results that validate our methods for quantizing the GS superstring; the agreement with gauge theory at one and two loops is impressive, though we will also confirm the previously observed breakdown of agreement at three-loop order.

In Sections 4.1 and 4.2 we present the details of the diagonalization of the perturbing string worldsheet Hamiltonian on degenerate subspaces of three-impurity states. We give a compressed discussion of general strategy, concentrating on the aspects of the problem which are new to the three-impurity case. An interesting new element is that the non-interacting degenerate subspace breaks up into several different supersymmetry multiplets so that the detailed accounting of multiplicities and irreducible decomposition amounts to a stringent test that the quantization has maintained the correct nonlinearly realized superconformal symmetries. Section 4.3 is devoted to the comparison of the string theory spectrum with gauge theory anomalous dimensions. We employ our own gauge theory data derived in Chapter 2 for the various higher-loop spin chains onto which the gauge theory anomalous dimension problem has been mapped. We find perfect agreement through two-loop order and, once again,

a breakdown at three loops.

Overall, the three-impurity regime of the string theory offers a much more stringent test of the duality away from the full plane-wave limit. While we are unable to offer (via this analysis) a solution to the disagreement with gauge theory at three loops, we can confirm that the complicated interacting worldsheet theory at $O(1/\hat{R}^2)$ in the curvature expansion is properly quantized and correct to two loops in λ .

4.1 Three-impurity spectrum: one loop in λ'

The three-impurity Fock space block-diagonalizes into separate spacetime fermion and spacetime boson sectors. The bosonic sector contains states that are purely bosonic (composed of three bosonic string oscillators) and states with bi-fermionic components:

$$a_q^{A\dagger} a_r^{B\dagger} a_s^{C\dagger} |J\rangle , \quad a_q^{A\dagger} b_r^{\alpha\dagger} b_s^{\beta\dagger} |J\rangle . \quad (4.1.1)$$

Pure boson states are mixed by the bosonic sector of the Hamiltonian H_{BB} , while states with bi-fermionic excitations are mixed both by the purely fermionic Hamiltonian H_{FF} and the bose-fermi sector H_{BF} . The sector of spacetime fermion states is composed of purely fermionic excitations and mixed states containing two bosonic oscillators:

$$b_q^{\alpha\dagger} b_r^{\beta\dagger} b_s^{\gamma\dagger} |J\rangle , \quad a_q^{A\dagger} a_r^{B\dagger} b_s^{\alpha\dagger} |J\rangle . \quad (4.1.2)$$

Pure fermion states are acted on by H_{FF} , and mixed states with bosonic excitations are acted on by H_{BB} and H_{BF} . This block diagonalization of the perturbing Hamiltonian is displayed schematically in table 4.1.

The three-impurity string states are subject to the usual level-matching condition on the mode indices: $q+r+s=0$. There are two generically different solutions of this constraint: all mode indices different ($q \neq r \neq s$) and two indices equal (e.g., $q=r=$

H_{int}	$a^{A\dagger}a^{B\dagger}a^{C\dagger} J\rangle$	$a^{A\dagger}b^{\alpha\dagger}b^{\beta\dagger} J\rangle$	$b^{\alpha\dagger}b^{\beta\dagger}b^{\gamma\dagger} J\rangle$	$a^{A\dagger}a^{B\dagger}b^{\alpha\dagger} J\rangle$
$\langle J a^Aa^Ba^C$	H_{BB}	H_{BF}	0	0
$\langle J a^Ab^\alpha b^\beta$	H_{BF}	$H_{\text{FF}} + H_{\text{BF}}$	0	0
$\langle J b^\alpha b^\beta b^\gamma$	0	0	H_{FF}	H_{BF}
$\langle J a^Aa^Bb^\alpha$	0	0	H_{BF}	$H_{\text{BB}} + H_{\text{BF}}$

Table 4.1: Three-impurity string states

$n, s = -2n$). In the inequivalent index case, there are $16^3 = 4,096$ degenerate states arising from different choices of spacetime labels on the mode creation operators. In the case of two equivalent indices, the dimension of the degenerate subspace is half as large (there are fewer permutations on mode indices that generate linearly independent states). The two types of basis break up into irreducible representations of $PSU(2,2|4)$ in different ways and must be studied separately.

As in the two-impurity case, the problem of diagonalizing the perturbation simplifies enormously when the matrix elements are expanded to leading order in λ' . We will take this approach here to obtain an overview of how degeneracies are lifted by the interaction. The generalization of the results to all loop orders in λ' (but still to first non-leading order in $1/J$) will be presented in the next section. It is once again the case that in the one-loop approximation, projection onto invariant subspaces under the manifest global $SO(4) \times SO(4)$ symmetry often diagonalizes the Hamiltonian directly (and at worst reduces it to a low-dimensional matrix). Symbolic manipulation programs were used to organize the complicated algebra and to perform explicit projections onto invariant subspaces.

4.1.1 Inequivalent mode indices ($q \neq r \neq s$)

In the sector of spacetime bosons, the subspace of purely bosonic states $a_q^{A\dagger}a_r^{B\dagger}a_s^{C\dagger}|J\rangle$ is 512-dimensional. When each of the three mode indices (q, r, s) are different, states with bi-fermionic excitations $a_q^{A\dagger}b_r^{\alpha\dagger}b_s^{\beta\dagger}|J\rangle$ are inequivalent under permutation of the mode indices, and form a 1,536-dimensional subsector. The entire bosonic

sector of the three-impurity state space therefore contains 2,048 linearly independent states. The fermionic sector decomposes in a similar manner: the subsector of purely fermionic states $b_q^{\alpha\dagger} b_r^{\beta\dagger} b_s^{\gamma\dagger} |J\rangle$ is 512-dimensional; fermionic states containing two bosonic excitations $a_q^{A\dagger} a_r^{B\dagger} b_s^{\alpha\dagger} |J\rangle$ are inequivalent under permutation of the mode indices, and comprise an additional 1,536-dimensional subsector. Adding this 2,048-dimensional fermion sector brings the dimensionality of the entire state space to 4,096.

Our first task is to evaluate the interaction Hamiltonian matrix. The matrix elements needed to fill out the spacetime boson sector are listed in table 4.2. To evaluate the entries, we express the Hamiltonian H_{int} computed in Chapter 3 in terms of mode creation and annihilation operators, expand the result in powers of λ' and then compute the indicated matrix elements between three-impurity Fock space states. The pure-boson (H_{BB}), pure-fermion (H_{FF}) and bose-fermi (H_{BF}) mixing sectors of H_{int} appear above in eqns. (3.4.25), (3.4.45) and (3.4.46) respectively. We collect below all the relevant results of this exercise.

H_{int}	$a_s^{D\dagger} a_r^{E\dagger} a_q^{F\dagger} J\rangle$	$a_s^{D\dagger} b_r^{\gamma\dagger} b_q^{\delta\dagger} J\rangle$	$a_r^{D\dagger} b_q^{\gamma\dagger} b_s^{\delta\dagger} J\rangle$	$a_r^{D\dagger} b_s^{\gamma\dagger} b_q^{\delta\dagger} J\rangle$
$\langle J a_q^A a_r^B a_s^C$	H_{BB}	H_{BF}	H_{BF}	H_{BF}
$\langle J a_q^A b_r^\alpha b_s^\beta$	H_{BF}	$H_{\text{FF}} + H_{\text{BF}}$	H_{BF}	H_{BF}
$\langle J a_s^A b_q^\alpha b_r^\beta$	H_{BF}	H_{BF}	$H_{\text{FF}} + H_{\text{BF}}$	H_{BF}
$\langle J a_r^A b_s^\alpha b_q^\beta$	H_{BF}	H_{BF}	H_{BF}	$H_{\text{FF}} + H_{\text{BF}}$

Table 4.2: H_{int} on spacetime-boson three-impurity string states ($q \neq r \neq s$)

We will use an obvious (m, n) matrix notation to distinguish the different entries

in table 4.2. The purely bosonic, 512-dimensional $(1, 1)$ block has the explicit form

$$\begin{aligned} \langle J | a_q^A a_r^B a_s^C (H_{\text{BB}}) a_s^{D\dagger} a_r^{E\dagger} a_q^{F\dagger} | J \rangle &= \frac{\lambda'}{J} \delta^{AF} \delta^{BE} \delta^{CD} (rs + q(r + s) - q^2 - r^2 - s^2) \\ &+ \frac{\lambda'}{2J} \left\{ \delta^{AF} \left[(r^2 + s^2) (\delta^{cd} \delta^{be} - \delta^{c'd'} \delta^{b'e'}) + (s^2 - r^2) (\delta^{be} \delta^{c'd'} - \delta^{cd} \delta^{b'e'}) \right. \right. \\ &\quad \left. \left. + 2rs (\delta^{bd} \delta^{ce} - \delta^{bc} \delta^{de} - \delta^{b'd'} \delta^{c'e'} + \delta^{b'c'} \delta^{d'e'}) \right] + (r \rightleftharpoons q, F \rightleftharpoons E, A \rightleftharpoons B) \\ &\quad \left. + (s \rightleftharpoons q, F \rightleftharpoons D, A \rightleftharpoons C) \right\}. \quad (4.1.3) \end{aligned}$$

The off-diagonal entries that mix purely bosonic states $a_q^{A\dagger} a_r^{B\dagger} a_s^{C\dagger} | J \rangle$ with states containing bi-fermions $a_q^{A\dagger} b_r^{\alpha\dagger} b_s^{\beta\dagger} | J \rangle$ are given by a separate set of 512-dimensional matrices. The $(1, 2)$ block in table 4.2, for example, yields

$$\langle J | a_q^A a_r^B a_s^C (H_{\text{BF}}) a_s^{D\dagger} b_r^{\alpha\dagger} b_q^{\beta\dagger} | J \rangle = \frac{\lambda'}{2J} \delta^{CD} q r \left\{ \left(\gamma^{ab'} \right)^{\alpha\beta} - \left(\gamma^{a'b} \right)^{\alpha\beta} \right\}, \quad (4.1.4)$$

where the index a (a') symbolizes the value of the vector index A , provided it is in the first (second) $SO(4)$. There are six blocks in this subsector, each given by a simple permutation of the mode indices (q, r, s) in eqn. (4.1.4). In table 4.2, these matrices occupy the $(1, 2)$, $(1, 3)$ and $(1, 4)$ blocks, along with their transposes in the $(2, 1)$, $(3, 1)$ and $(4, 1)$ entries.

The pure-fermion sector of the Hamiltonian, H_{FF} , has non-vanishing matrix elements between states containing bi-fermionic excitations. The H_{FF} contribution to the $(2, 2)$ block, for example, is given by

$$\begin{aligned} \langle J | b_q^\alpha b_r^\beta a_s^A (H_{\text{FF}}) a_s^{B\dagger} b_r^{\gamma\dagger} b_q^{\delta\dagger} | J \rangle &= -\frac{\lambda'}{2J} (q - r)^2 \delta^{AB} \delta^{\alpha\delta} \delta^{\gamma\beta} \\ &+ \frac{\lambda'}{24J} \delta^{AB} q r \left\{ (\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\beta\delta} - (\gamma^{ij})^{\alpha\beta} (\gamma^{ij})^{\gamma\delta} - (\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\beta\gamma} \right. \\ &\quad \left. - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\beta\delta} + (\gamma^{i'j'})^{\alpha\beta} (\gamma^{i'j'})^{\gamma\delta} + (\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\beta\gamma} \right\}. \quad (4.1.5) \end{aligned}$$

A similar contribution, related to this one by simple permutations of the mode indices (q, r, s) , appears in the diagonal blocks $(3, 3)$ and $(4, 4)$ as well.

The bose-fermi mixing Hamiltonian H_{BF} makes the following contribution to the lower diagonal blocks (2, 2), (3, 3) and (4, 4) in table 4.2:

$$\begin{aligned} \langle J | b_q^\alpha b_r^\beta a_s^A (H_{\text{BF}}) a_s^{B\dagger} b_r^{\gamma\dagger} b_q^{\delta\dagger} | J \rangle = & \frac{\lambda'}{2J} \left\{ 2s(q+r-s) \delta^{ab} \delta^{\alpha\delta} \delta^{\beta\gamma} \right. \\ & - r s \left[(\gamma^{ab})^{\beta\gamma} - (\gamma^{a'b'})^{\beta\gamma} \right] - s q \left[(\gamma^{ab})^{\alpha\delta} - (\gamma^{a'b'})^{\alpha\delta} \right] \\ & \left. - 2 \left[q^2 + r^2 + s^2 - s(q+r) \right] \delta^{a'b'} \delta^{\alpha\delta} \delta^{\beta\gamma} \right\}. \end{aligned} \quad (4.1.6)$$

The H_{BF} sector also makes the following contribution to the off-diagonal (2, 3) block:

$$\langle J | b_q^\alpha b_r^\beta a_s^A (H_{\text{BF}}) a_r^{B\dagger} b_q^{\gamma\dagger} b_s^{\delta\dagger} | J \rangle = -\frac{\lambda'}{2J} \delta^{\alpha\gamma} r s \left\{ (\delta^{ab} - \delta^{a'b'}) \delta^{\beta\delta} - (\gamma^{ab})^{\beta\delta} + (\gamma^{a'b'})^{\beta\delta} \right\}. \quad (4.1.7)$$

The contributions of H_{BF} to the remaining off-diagonal blocks (2, 3), (2, 4), etc. are obtained by appropriate index permutations.

The sector of spacetime fermions decomposes in a similar fashion. The fermion analogue of table 4.2 for the bosonic sector appears in table 4.3. The (1, 1) fermion

H_{int}	$b_s^{\zeta\dagger} b_r^{\epsilon\dagger} b_q^{\delta\dagger} J \rangle$	$a_s^{C\dagger} a_r^{D\dagger} b_q^{\delta\dagger} J \rangle$	$a_r^{C\dagger} a_q^{D\dagger} b_s^{\delta\dagger} J \rangle$	$a_r^{C\dagger} a_s^{D\dagger} b_q^{\delta\dagger} J \rangle$
$\langle J b_q^\alpha b_r^\beta b_s^\gamma$	H_{FF}	H_{BF}	H_{BF}	H_{BF}
$\langle J b_q^\alpha a_r^A a_s^B$	H_{BF}	$H_{\text{BB}} + H_{\text{BF}}$	H_{BF}	H_{BF}
$\langle J b_s^\alpha a_q^A a_r^B$	H_{BF}	H_{BF}	$H_{\text{BB}} + H_{\text{BF}}$	H_{BF}
$\langle J b_r^\alpha a_s^A a_q^B$	H_{BF}	H_{BF}	H_{BF}	$H_{\text{BB}} + H_{\text{BF}}$

Table 4.3: Interaction Hamiltonian on spacetime fermion three-impurity states ($q \neq r \neq s$)

block is occupied by the pure-fermion sector of the Hamiltonian taken between the

purely fermionic three-impurity states $b_q^{\alpha\dagger} b_r^{\beta\dagger} b_s^{\gamma\dagger} |J\rangle$:

$$\begin{aligned} \langle J | b_q^\alpha b_r^\beta b_s^\gamma (H_{\text{FF}}) b_s^{\zeta\dagger} b_r^{\epsilon\dagger} b_q^{\delta\dagger} | J \rangle &= -\frac{\lambda'}{J} [q^2 + r^2 + s^2 - rs - q(r+s)] \delta^{\alpha\delta} \delta^{\beta\epsilon} \delta^{\gamma\zeta} \\ &+ \frac{\lambda'}{24J} \delta^{\alpha\delta} rs \left\{ (\gamma^{ij})^{\beta\gamma} (\gamma^{ij})^{\epsilon\zeta} - (\gamma^{ij})^{\beta\epsilon} (\gamma^{ij})^{\gamma\zeta} + (\gamma^{ij})^{\beta\zeta} (\gamma^{ij})^{\gamma\epsilon} \right. \\ &\quad - (\gamma^{i'j'})^{\beta\gamma} (\gamma^{i'j'})^{\epsilon\zeta} + (\gamma^{i'j'})^{\beta\epsilon} (\gamma^{i'j'})^{\gamma\zeta} - (\gamma^{i'j'})^{\beta\zeta} (\gamma^{i'j'})^{\gamma\epsilon} \\ &\quad \left. + (r \rightleftharpoons q, \alpha \rightleftharpoons \beta, \delta \rightleftharpoons \epsilon) + (s \rightleftharpoons q, \alpha \rightleftharpoons \gamma, \delta \rightleftharpoons \zeta) \right\}. \end{aligned} \quad (4.1.8)$$

The off-diagonal (1, 2), (1, 3) and (1, 4) blocks (and their transposes) mix purely fermionic states with $a_s^{A\dagger} a_r^{B\dagger} b_q^{\alpha\dagger} |J\rangle$ states:

$$\langle J | b_q^\alpha b_r^\beta b_s^\gamma (H_{\text{BF}}) a_s^{A\dagger} a_r^{B\dagger} b_q^{\alpha\dagger} | J \rangle = -\frac{\lambda'}{2J} \delta^{\alpha\delta} rs \left\{ (\gamma^{ab'})^{\beta\gamma} - (\gamma^{a'b})^{\beta\gamma} \right\}. \quad (4.1.9)$$

The lower-diagonal (2, 2), (3, 3) and (4, 4) blocks receive contributions from the pure boson sector of the Hamiltonian:

$$\begin{aligned} \langle J | b_q^\alpha a_r^A a_s^B (H_{\text{BB}}) a_s^{C\dagger} a_r^{D\dagger} b_q^{\beta\dagger} | J \rangle &= -\frac{\lambda'}{2J} \delta^{\alpha\beta} \left\{ (r-s)^2 \delta^{BC} \delta^{AD} \right. \\ &\quad - (r^2 + s^2) (\delta^{ad} \delta^{bc} - \delta^{a'd'} \delta^{b'c'}) \\ &\quad - 2rs (\delta^{ac} \delta^{bd} - \delta^{ab} \delta^{cd} - \delta^{a'c'} \delta^{b'd'} + \delta^{a'b'} \delta^{c'd'}) \\ &\quad \left. + (r^2 - s^2) (\delta^{ad} \delta^{b'c'} - \delta^{a'd'} \delta^{bc}) \right\}. \end{aligned} \quad (4.1.10)$$

In the same diagonal blocks of table 4.3, the H_{BF} sector contributes

$$\begin{aligned} \langle J | b_q^\alpha a_r^A a_s^B (H_{\text{BF}}) a_s^{C\dagger} a_r^{D\dagger} b_q^{\beta\dagger} | J \rangle &= \frac{\lambda'}{8J} \left\{ \delta^{\alpha\beta} \left[+ (8q(r+s) - 5(r^2 + s^2) - 6q^2) \delta^{AD} \delta^{BC} \right. \right. \\ &\quad + (3q^2 + s^2) \delta^{AD} \delta^{bc} + (3q^2 + r^2) \delta^{BC} \delta^{ad} + (r^2 - 5q^2) \delta^{BC} \delta^{a'd'} \\ &\quad \left. + (s^2 - 5q^2) \delta^{AD} \delta^{b'c'} \right] - 4\delta^{BC} qr \left[(\gamma^{ad})^{\alpha\beta} - (\gamma^{a'd'})^{\alpha\beta} \right] \\ &\quad \left. - 4\delta^{AD} qs \left[(\gamma^{bc})^{\alpha\beta} - (\gamma^{b'c'})^{\alpha\beta} \right] \right\}. \end{aligned} \quad (4.1.11)$$

Finally, the off-diagonal blocks (2, 3), (2, 4) and (3, 4) (plus their transpose entries) are given by the H_{BF} matrix element

$$\begin{aligned} \langle J | b_q^\alpha a_r^A a_s^B (H_{\text{BF}}) a_r^{C\dagger} a_q^{D\dagger} b_s^{\beta\dagger} | J \rangle &= -\frac{\lambda'}{32J} \delta^{AC} \left\{ \delta^{\alpha\beta} \left[(q-s)^2 \delta^{BD} - (q^2 + 14qs + s^2) \delta^{bd} \right. \right. \\ &\quad \left. \left. - (q^2 - 18qs + s^2) \delta^{b'd'} \right] + 16qs \left[(\gamma^{bd})^{\alpha\beta} - (\gamma^{b'd'})^{\alpha\beta} \right] \right\}. \end{aligned} \quad (4.1.12)$$

A significant departure from the two-impurity case is that all these matrix elements have, along with their spacetime index structures, nontrivial dependence on the mode indices. The eigenvalues could potentially have very complicated mode-index dependence but, as we shall see, they do not. This amounts to a rigid consistency check on the whole procedure that was not present in the two-impurity case.

4.1.2 Matrix diagonalization: inequivalent modes ($q \neq r \neq s$)

We now turn to the task of diagonalizing the one-loop approximation to the perturbing Hamiltonian. To simplify the task, we exploit certain block diagonalizations that hold to leading order in λ' (but not to higher orders). While we eventually want to study the spectrum to all orders in λ' , diagonalizing the Hamiltonian at one loop will reveal the underlying supermultiplet structure. As an example of the simplifications we have in mind, we infer from (4.1.4) that the matrix elements of H_{BF} between pure boson states $a_q^{A\dagger} a_r^{B\dagger} a_s^{C\dagger} | J \rangle$ and bifermionic spacetime bosons vanish to leading order in λ' if all three $SO(8)$ bosonic vector indices lie within the same $SO(4)$, descended either from AdS_5 or S^5 . Restricting to such states brings the bosonic sector of the Hamiltonian into the block-diagonal form in table 4.4. This leaves two 64-dimensional subspaces of purely bosonic states on which the perturbation is block diagonal, as recorded in table 4.5.

H_{int}	$a^{a\dagger}a^{b\dagger}a^{c\dagger} J\rangle + a^{a'\dagger}a^{b'\dagger}a^{c'\dagger} J\rangle$	$a^{A\dagger}b^{\alpha\dagger}b^{\beta\dagger} J\rangle$
$\langle J a^a a^b a^c + \langle J a^{a'} a^{b'} a^{c'}$	H_{BB}	0
$\langle J a^A b^\alpha b^\beta$	0	$H_{\text{FF}} + H_{\text{BF}}$

Table 4.4: Block-diagonal $SO(4)$ projection on bosonic three-impurity string states

H_{int}	$a^{a\dagger}a^{b\dagger}a^{c\dagger} J\rangle$	$a^{a'\dagger}a^{b'\dagger}a^{c'\dagger} J\rangle$
$\langle J a^a a^b a^c$	$(H_{\text{BB}})_{64 \times 64}$	0
$\langle J a^{a'} a^{b'} a^{c'}$	0	$(H_{\text{BB}})_{64 \times 64}$

Table 4.5: $SO(4)$ projection on purely bosonic states

Since the interaction Hamiltonian has manifest $SO(4) \times SO(4)$ symmetry, it is useful to project matrix elements onto irreps of that group before diagonalizing. In some cases the irrep is unique, and projection directly identifies the corresponding eigenvalue. In the cases where an irrep has multiple occurrences, there emerges an unavoidable matrix diagonalization that is typically of low dimension. In what follows, we will collect the results of carrying out this program on the one-loop interaction Hamiltonian. A very important feature of the results that appear is that all the eigenvalues turn out to have a common simple dependence on mode indices. More precisely, the expansion of the eigenvalues for inequivalent mode indices (q, r, s) out to first non-leading order in λ' and $1/J$ can be written as

$$E_J(q, r, s) = 3 + \frac{\lambda'(q^2 + r^2 + s^2)}{2} \left(1 + \frac{\Lambda}{J} + O(J^{-2}) \right), \quad (4.1.13)$$

where, as in Chapter 3, Λ is a pure number that characterizes the lifting of the degeneracy in the various sectors. (The notation Λ_{BB} , Λ_{BF} and Λ_{FF} will again be used to denote energy corrections arising entirely from the indicated sectors of the perturbing Hamiltonian.) This simple quadratic dependence of the eigenvalues on the mode indices does not automatically follow from the structure of the matrix elements themselves, but is important for the successful match to gauge theory eigenvalues. In

what follows, we will catalog some of the different Λ values that occur, along with their $SO(4) \times SO(4)$ irreps (and multiplicities). When we have the complete list, we will discuss how they are organized into supermultiplets.

In the $SO(4)$ projection in table 4.5, we will find a set of 64 eigenvalues for both the $SO(4)_{AdS}$ and $SO(4)_{S^5}$ subsectors. We record this eigenvalue spectrum in table 4.6, using a now-familiar $SU(2)^2 \times SU(2)^2$ notation. For comparison, it is displayed alongside the projection of the two-impurity spectrum onto the same subspace (as found in Chapter 3). In the three-impurity case, the $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ level in

$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{BB}	$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{BB}
$(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$	-8	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-6
$[\mathbf{1}, \mathbf{1}; (\mathbf{2} + \mathbf{4}), \mathbf{2}] + [\mathbf{1}, \mathbf{1}; \mathbf{2}, (\mathbf{2} + \mathbf{4})]$	-6	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	-4
$[\mathbf{1}, \mathbf{1}; (\mathbf{2} + \mathbf{4}), (\mathbf{2} + \mathbf{4})]$	-4	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$	-2
$[(\mathbf{2} + \mathbf{4}), (\mathbf{2} + \mathbf{4}); \mathbf{1}, \mathbf{1}]$	-2	$(\mathbf{3}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-2
$[(\mathbf{2} + \mathbf{4}), \mathbf{2}; \mathbf{1}, \mathbf{1}] + [\mathbf{2}, (\mathbf{2} + \mathbf{4}); \mathbf{1}, \mathbf{1}]$	0	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	0
$(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$	2	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	2

Table 4.6: Three-impurity energy spectrum in the pure-boson $SO(4)$ projection (left panel) and two-impurity energy spectrum in the same projection (right panel)

the $SO(4)_{S^5}$ subsector clearly descends from the two-impurity singlet $(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$ in the same $SO(4)$ subgroup. In the same manner, the three-impurity $[\mathbf{1}, \mathbf{1}; (\mathbf{2} + \mathbf{4}), \mathbf{2}] + [\mathbf{1}, \mathbf{1}; \mathbf{2}, (\mathbf{2} + \mathbf{4})]$ level descends from the $SO(4)_{S^5}$ antisymmetric two-impurity state $(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$, and the three-impurity $[\mathbf{1}, \mathbf{1}; (\mathbf{2} + \mathbf{4}), (\mathbf{2} + \mathbf{4})]$ level is tied to the two-impurity symmetric-traceless $(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$ irrep. In the $SO(4)_{S^5}$ subsector, each of these levels receives a shift to the energy of -2 . The total multiplicity of each of these levels is also increased by a factor of four when the additional $(\mathbf{2}, \mathbf{2})$ is tensored into the two-impurity state space. The $SO(4)_{AdS}$ subsector follows a similar pattern: the $(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$, $[(\mathbf{2} + \mathbf{4}), \mathbf{2}; \mathbf{1}, \mathbf{1}] + [\mathbf{2}, (\mathbf{2} + \mathbf{4}); \mathbf{1}, \mathbf{1}]$ and $[(\mathbf{2} + \mathbf{4}), (\mathbf{2} + \mathbf{4}); \mathbf{1}, \mathbf{1}]$ levels appear as three-impurity descendants of the two-impurity irrep spectrum $(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1}) + (\mathbf{3}, \mathbf{3}; \mathbf{1}, \mathbf{1})$. In this subsector, however, the three-impurity

energies are identical to those in the two-impurity theory.

The bosonic $SO(4)$ projection has a precise fermionic analogue. Similar to the bosons, the $SO(9, 1)$ spinors b_q^\dagger decompose as $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) + (\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ under the action of Π parity:

$$\Pi \hat{b}_q^\dagger = \hat{b}_q^\dagger, \quad \Pi \tilde{b}_q^\dagger = -\tilde{b}_q^\dagger. \quad (4.1.14)$$

(As described above, the notation \hat{b}_q^\dagger labels $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ spinors with positive eigenvalue under Π , and \tilde{b}_q^\dagger indicates $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ spinors, which are negative under Π .) Analogous to the $SO(4)$ projection on the $SO(8)$ bosonic operators $a_q^{A\dagger} \rightarrow a_q^{a\dagger} + a_q^{a'\dagger}$, projecting out the positive or negative eigenvalues of Π on the eight-component spinor $b_q^{\alpha\dagger}$ leaves a subspace of four-component spinors spanned by \hat{b}_q^\dagger and \tilde{b}_q^\dagger .

We can perform a projection on the subsector in table 4.2 similar to that appearing in table 4.5. In this case, instead of three bosonic impurities mixing with a single bosonic (plus a bi-fermionic) excitation, we are now interested in projecting out particular interactions between a purely fermionic state and a state with one fermionic and two bosonic excitations. Using \pm to denote the particular representation of the fermionic excitations, the off-diagonal elements given by (4.1.9) vanish for $+++ \rightarrow \pm$ and $--- \rightarrow \pm$ interactions. In other words, the pure fermion states in the $(1, 1)$ block of table 4.3 will not mix with states containing two bosonic excitations if all three fermionic oscillators lie in the same Π projection. This projection appears schematically in table 4.7.

H_{int}	$\hat{b}^{\alpha\dagger}\hat{b}^{\beta\dagger}\hat{b}^{\gamma\dagger} J\rangle + \tilde{b}^{\alpha\dagger}\tilde{b}^{\beta\dagger}\tilde{b}^{\gamma\dagger} J\rangle$	$a^{A\dagger}a^{B\dagger}b^{\alpha\dagger} J\rangle$
$\langle J \hat{b}^\alpha\hat{b}^\beta\hat{b}^\gamma + \langle J \tilde{b}^\alpha\tilde{b}^\beta\tilde{b}^\gamma$	H_{FF}	0
$\langle J a^A a^B b^\alpha$	0	$H_{\text{BB}} + H_{\text{BF}}$

Table 4.7: Block-diagonal projection on fermionic three-impurity string states

The $(1, 1)$ pure fermion block in table 4.7 breaks into two 64-dimensional subsectors under this projection. By tensoring an additional $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ or $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$

impurity into the two-impurity state space, we expect to see a multiplicity structure in this projection given by

$$\begin{aligned}
(\mathbf{1}, \mathbf{2}) \times (\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2}) &= (\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2}) + [\mathbf{1}, \mathbf{2}; \mathbf{1}, (\mathbf{2} + \mathbf{4})] \\
&\quad + [\mathbf{1}, (\mathbf{2} + \mathbf{4}); \mathbf{1}, \mathbf{2}] + [\mathbf{1}, (\mathbf{2} + \mathbf{4}); \mathbf{1}, (\mathbf{2} + \mathbf{4})] , \\
(\mathbf{2}, \mathbf{1}) \times (\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) &= (\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) + [\mathbf{2}, \mathbf{1}; (\mathbf{2} + \mathbf{4}), \mathbf{1}] \\
&\quad + [(\mathbf{2} + \mathbf{4}), \mathbf{1}; \mathbf{2}, \mathbf{1}] + [(\mathbf{2} + \mathbf{4}), \mathbf{1}; (\mathbf{2} + \mathbf{4}), \mathbf{1}] , \quad (4.1.15)
\end{aligned}$$

for a total of 128 states. The projections onto the two 64-dimensional Π_+ and Π_- subspaces yield identical eigenvalues and multiplicities. The results for both subspaces are presented in table 4.8: The two-impurity bi-fermion states in table 4.8 are

$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{FF}	$SO(4)_{AdS} \times SO(4)_{S^5}$	Λ_{FF}
$(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1}) + (\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$	-3	$(\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{1})$	-2
$[\mathbf{2}, \mathbf{1}; (\mathbf{2} + \mathbf{4}), \mathbf{1}] + [\mathbf{1}, \mathbf{2}; \mathbf{1}, (\mathbf{2} + \mathbf{4})]$	-1	$(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{1}; \mathbf{1}, \mathbf{3})$	0
$[(\mathbf{2} + \mathbf{4}), \mathbf{1}; \mathbf{2}, \mathbf{1}] + [\mathbf{1}, (\mathbf{2} + \mathbf{4}); \mathbf{1}, \mathbf{2}]$	-5	$(\mathbf{3}, \mathbf{1}; \mathbf{1}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{1})$	-4
$[(\mathbf{2} + \mathbf{4}), \mathbf{1}; (\mathbf{2} + \mathbf{4}), \mathbf{1}] + [\mathbf{1}, (\mathbf{2} + \mathbf{4}); \mathbf{1}, (\mathbf{2} + \mathbf{4})]$	-3	$(\mathbf{3}, \mathbf{1}; \mathbf{3}, \mathbf{1}) + (\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{3})$	-2

Table 4.8: Spectrum of three-impurity states (left panel) and two-impurity states (right panel) created by Π_{\pm} -projected fermionic creation operators

spacetime bosons while the tri-fermion states are spacetime fermions. For comparison purposes, we have displayed both spectra. Note that the $O(1/J)$ energy corrections of the two types of state are simply displaced by -1 relative to each other.

This exhausts the subspaces that can be diagonalized by simple irrep projections. The remaining eigenvalues must be obtained by explicit diagonalization of finite dimensional submatrices obtained by projection onto representations with multiple occurrence. The upshot of these more complicated eigenvalue calculations is that the first-order λ' eigenvalues take on all integer values from $\Lambda = -8$ to $\Lambda = +2$, alternating between spacetime bosons and fermions as Λ is successively incremented by one unit.

4.1.3 Assembling eigenvalues into supermultiplets

Finally, we need to understand how the perturbed three-impurity spectrum breaks up into extended supersymmetry multiplets. This is relatively easy to infer from the multiplicities of the perturbed eigenvalues (and the multiplicities are a side result of the calculation of the eigenvalues themselves). In the last subsection, we described a procedure for diagonalizing the one-loop perturbing Hamiltonian on the 4,096-dimensional space of three-impurity string states with mode indices $p \neq q \neq r$. The complete results for the eigenvalues Λ and their multiplicities are stated in table 4.9 (we use the notation of (4.1.13), while the B and F subscripts are used to indicate bosonic and fermionic levels in the supermultiplet).

Λ	-8	-7	-6	-5	-4	-3	-2	-1	0	1	2
Multiplicity	4_B	40_F	180_B	480_F	840_B	1008_F	840_B	480_F	180_B	40_F	4_B

Table 4.9: Complete three-impurity energy spectrum (with multiplicities)

The Λ eigenvalues in table 4.9 are integer-spaced, which is consistent with supersymmetry requirements (for details, the reader is referred to the discussion following eqn. (3.5.16) in Chapter 3 above). However, because the range between top and bottom eigenvalues is ten, rather than eight, the 4,096-dimensional space must be built on more than one type of extended supermultiplet, with more than one choice of c in the general formula $\Lambda = L + c$ (where L is the supermultiplet level and c is some numerical constant). This is to be contrasted with the two-impurity case, where the degenerate space was exactly 256-dimensional and was spanned by a single superconformal primary whose lowest member was a singlet under both Lorentz transformations and the residual $SO(4)$ R -symmetry. We can readily infer what superconformal primaries are needed to span the degenerate three-impurity state space by applying a little numerology to table 4.9. The lowest eigenvalue is $\Lambda = -8$: it has multiplicity 4 and, according to table 4.6, its $SO(4) \times SO(4)$ decomposition is $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ (spacetime scalar, R -charge $SO(4)$ four-vector). According to the gen-

eral arguments about how the full extended supermultiplet is built by acting on a “bottom” state with the eight raising operators, it is the base of a supermultiplet of 4×256 states extending up to $\Lambda = 0$. By the same token, there is a highest eigenvalue $\Lambda = +2$: it has multiplicity 4 and, according to table 4.6, its $SO(4) \times SO(4)$ decomposition is $(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$ (spacetime vector, R -charge singlet). Using lowering operators instead of raising operators, we see that one derives from it a supermultiplet of 4×256 operators with eigenvalues extending from $\Lambda = -6$ to $\Lambda = +2$. The multiplicities of the Λ eigenvalues occurring in these two supermultiplets are of course given by binomial coefficients, as described above. By comparing with the *total* multiplicities of each allowed Λ (as listed in table 4.9) we readily see that what remains are 8×256 states with eigenvalues running from $\Lambda = -7$ to $\Lambda = +1$ with the correct binomial coefficient pattern of multiplicities. The top and bottom states here are spacetime fermions and must lie in a spinor representation of the Lorentz group. It is not hard to see that they lie in the eight-dimensional $SO(4) \times SO(4)$ irrep $(\mathbf{2}, \mathbf{1}; \mathbf{1}, \mathbf{2}) + (\mathbf{1}, \mathbf{2}; \mathbf{2}, \mathbf{1})$. This exhausts all the states and we conclude that the three-impurity state space is spanned by three distinct extended superconformal multiplets. The detailed spectrum is given in table 4.10 (where the last line records the total multiplicity at each level as given in table 4.9 and the first line records the two-impurity spectrum for reference). Note the peculiar feature that certain energies are shared by all three multiplets: this is an accidental degeneracy that does not survive at higher loop order.

Λ	-8	-7	-6	-5	-4	-3	-2	-1	0	1	2	
$\Delta_0 = 2$			1_B	8_F	28_B	56_F	70_B	56_F	28_B	8_F	1_B	scalar
$\Delta_0 = 3$	4	32	112	224	280	224	112	32	4			$SO(4)_{S^5}$ vector
			4	32	112	224	280	224	112	32	4	$SO(4)_{AdS_5}$ vector
			8	64	224	448	560	448	224	64	8	spinor
Total	4_B	40_F	180_B	480_F	840_B	1008_F	840_B	480_F	180_B	40_F	4_B	4,096

Table 4.10: Submultiplet breakup of the three-impurity spectrum

A complete analysis of the agreement with gauge theory anomalous dimensions will have to be deferred until a later section: the dimensions of three-impurity gauge

theory operators are much harder to calculate than those of the two-impurity operators and there are few results in the literature, even at one loop. However, it is worth making a few preliminary remarks at this point. Since there are three superconformal multiplets, we have only three independent anomalous dimensions to compute. Minahan and Zarembo [39] found that the problem simplifies dramatically if we study the one-loop anomalous dimension of the special subset of single-trace operators of the form $\text{tr}(\phi^I Z^J)$ (and all possible permutations of the fields inside the trace), where the R -charge is carried by an $SO(4) \times SO(4)$ singlet scalar field Z and the impurities are insertions of a scalar field ϕ lying in the $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ (vector) irrep of the residual $SO(4) \times SO(4)$ symmetry. More formally, these operators are in the $SO(4) \times SO(4)$ irrep obtained by completely symmetrizing I vectors in the $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ irrep. The crucial point is that such operators form a closed sector, mixing only among themselves under the anomalous dimension operator. More importantly, the action of the one-loop anomalous dimension operator on this closed sector can be recast as the action of an integrable spin chain Hamiltonian of a type solvable by Bethe ansatz techniques. Although the Bethe ansatz is generally not analytically soluble, Minahan and Zarembo used it to obtain a virial expansion for the anomalous dimension in which the number I of impurities is held fixed, while the R -charge J is taken to be large (see eqn. (5.29) in [39]). In terms of the number of spin chain lattice sites K , their result appears as

$$\gamma_{\text{so}(6)} = \frac{\lambda}{2K^3} \sum_n M_n k_n^2 (K + M_n + 1) + O(K^{-4}) . \quad (4.1.16)$$

The integer k_n represents pseudoparticle momenta on the spin chain, and is dual to the string theory worldsheet mode indices; the quantity M_n labels the number of trace impurities with identical k_n . With I impurities, the spin chain length is given in terms of the R -charge by $K = J + I$, which leads to

$$\gamma_{\text{so}(6)} = \frac{\lambda}{2J^3} \sum_n M_n k_n^2 (J - 2I + M_n + 1) + O(J^{-4}) . \quad (4.1.17)$$

This virial expansion is similar in character to (4.1.13) and, for $I = 3$ (the three-impurity case), it matches that equation precisely with $\Lambda = -4$.

On the string theory side, three completely symmetrized $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ vectors form a tensor in the $(\mathbf{1}, \mathbf{1}; \mathbf{4}, \mathbf{4})$ irrep; such an irrep can be constructed from three $SO(4)_{S^5}$ vector (bosonic) creation operators. Table 4.6 shows that the corresponding string perturbation theory eigenvalue is (at one-loop order) $\Lambda = -4$ as well. We infer from table 4.10 that this eigenvalue lies at level $L = 4$ of the $SO(4)_{S^5}$ vector superconformal multiplet (and this argument takes care of the gauge theory/string theory comparison for all other operators in that multiplet).

The sector described above is often called an $\mathfrak{so}(6)^2$ sector on the gauge theory side, with reference to the subalgebra of the full superconformal algebra under which it is invariant. In an $\mathfrak{su}(2)$ subspace of the $\mathfrak{so}(6)$, this sector becomes closed to all loop order. For future reference, we note that Beisert [69] has identified two other closed sectors of operators in the gauge theory. In addition to the bosonic $\mathfrak{su}(2)$ sector, a bosonic $\mathfrak{sl}(2)$ sector and an $\mathfrak{su}(2|3)$ sector (of which the closed $\mathfrak{su}(2)$ sector is a subsector) are also exactly closed. It should be noted that integrable $\mathfrak{sl}(2)$ spin chains were discovered some time ago in phenomenologically motivated studies of the scaling behavior of high-energy scattering amplitudes in physical, non-supersymmetric QCD [90] (see also [91–94]). The $\mathfrak{su}(2|3)$ spin chain was studied more recently in [33]: this closed sector breaks into the $\mathfrak{su}(2)$ bosonic sector and a special fermionic subsector, which we denote as $\mathfrak{su}(1|1)$ (a subalgebra of $\mathfrak{su}(2|3)$).

In the string theory, the subsectors analogous to the gauge theory $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ are constructed out of completely symmetrized $SO(4)_{AdS}$ bosons and completely symmetrized fermions of the same Π eigenvalue, respectively (see Chapter 3 or ref. [27]). They correspond to the central $L = 4$ levels of the remaining two supermultiplets in table 4.10, and a calculation of their eigenvalues would complete the analysis of the match between three-impurity operators and string states at one-loop order. Since

²This notation is used to distinguish the protected gauge theory symmetry groups from those in the string theory.

we eventually want to go beyond one loop, where Bethe ansatz technology is less well-developed, we have found it useful to employ the numerical methods presented in Chapter 2 for evaluating spin chain eigenvalues (we refer the reader to Chapter 2, or to ref. [95], for a check of our results against Bethe-ansatz techniques, including the higher-loop corrections of [34]). This subject will be developed in a later section.

4.1.4 Two equivalent mode indices ($q = r = n$, $s = -2n$)

When two mode indices are allowed to be equal, the analysis becomes slightly more complicated. Since we are diagonalizing a Hamiltonian that is quartic in oscillators in a basis of three-impurity string states, one oscillator in the “in” state must always be directly contracted with one oscillator in the “out” state and, with two equal mode indices, there are many more nonvanishing contributions to each matrix element. While the matrix elements are more complicated, the state space is only half as large when two mode indices are allowed to be equal (only half as many mode-index permutations on the basis states generate linearly independent states). As a result, the fermionic and bosonic sectors of the Hamiltonian are each 1,024-dimensional. By the same token, the multiplet structure of the energy eigenstates will be significantly different from the unequal mode index case studied in the previous subsection.

To study this case, we make the mode index choice

$$q = r = n , \quad s = -2n . \quad (4.1.18)$$

The structure of matrix elements of the string Hamiltonian between spacetime bosons is given in table 4.11. This table seems to describe a 3×3 block matrix with 512×512 blocks in each subsector, giving a 1,536-dimensional state space. However, the vector and spinor indices are required to run over values that generate linearly independent basis states. This eliminates one third of the possible index assignments, implying that the matrix is in fact $1,024 \times 1,024$.

To evaluate the entries in table 4.11, we express the Hamiltonians (3.4.25, 3.4.45,

H_{int}	$a_{-2n}^{D\dagger} a_n^{E\dagger} a_n^{F\dagger} J\rangle$	$a_{-2n}^{D\dagger} b_n^{\gamma\dagger} b_n^{\delta\dagger} J\rangle$	$a_n^{D\dagger} b_n^{\gamma\dagger} b_{-2n}^{\delta\dagger} J\rangle$
$\langle J a_n^A a_n^B a_{-2n}^C$	H_{BB}	H_{BF}	H_{BF}
$\langle J a_n^A b_n^\alpha b_{-2n}^\beta$	H_{BF}	$H_{\text{FF}} + H_{\text{BF}}$	H_{BF}
$\langle J a_{-2n}^A b_n^\alpha b_n^\beta$	H_{BF}	H_{BF}	$H_{\text{FF}} + H_{\text{BF}}$

Table 4.11: Bosonic three-impurity string perturbation matrix with ($q = r = n$, $s = -2n$)

3.4.46) in terms of mode creation and annihilation operators, expand the result in powers of λ' and compute the indicated matrix elements between three-impurity Fock space states. We collect below all the relevant results of this exercise for this equal-mode-index case.

The purely bosonic subsector in the (1, 1) block is given by

$$\begin{aligned}
\langle J | a_n^A a_n^B a_{-2n}^C (H_{\text{BB}}) a_{-2n}^{D\dagger} a_n^{E\dagger} a_n^{F\dagger} | J \rangle = & \frac{n^2 \lambda}{2J} \left\{ 5 \delta^{BF} \delta^{cd} \delta^{ae} + 5 \delta^{AF} \delta^{cd} \delta^{be} - 4 \delta^{BF} \delta^{ad} \delta^{ce} \right. \\
& + 4 \delta^{BF} \delta^{ac} \delta^{de} + 4 \delta^{AF} \delta^{bc} \delta^{de} + 5 \delta^{BE} \delta^{cd} \delta^{af} - 4 \delta^{BE} \delta^{ad} \delta^{cf} + 4 \delta^{BE} \delta^{ac} \delta^{df} \\
& + 4 \delta^{AE} \delta^{bc} \delta^{df} - 4 \delta^{bd} \left(\delta^{AF} \delta^{ce} + \delta^{AE} \delta^{cf} \right) + 3 \delta^{BF} \delta^{ae} \delta^{c'd'} + 3 \delta^{AF} \delta^{be} \delta^{c'd'} \\
& + 3 \delta^{BE} \delta^{af} \delta^{c'd'} - 3 \delta^{BF} \delta^{cd} \delta^{a'e'} - 3 \delta^{AF} \delta^{cd} \delta^{b'e'} - 5 \delta^{BF} \delta^{c'd'} \delta^{a'e'} - 5 \delta^{AF} \delta^{c'd'} \delta^{b'e'} \\
& + 4 \delta^{BF} \delta^{a'd'} \delta^{c'e'} + 4 \delta^{AF} \delta^{b'd'} \delta^{c'e'} - 4 \delta^{BF} \delta^{a'c'} \delta^{d'e'} - 4 \delta^{AF} \delta^{b'c'} \delta^{d'e'} - 3 \delta^{BE} \delta^{cd} \delta^{a'f'} \\
& - 3 \delta^{AE} \delta^{cd} \delta^{b'f'} - 5 \delta^{BE} \delta^{c'd'} \delta^{a'f'} - 5 \delta^{AE} \delta^{c'd'} \delta^{b'f'} + 4 \delta^{BE} \delta^{a'd'} \delta^{c'f'} + 4 \delta^{AE} \delta^{b'd'} \delta^{c'f'} \\
& - 4 \delta^{BE} \delta^{a'c'} \delta^{d'f'} - 4 \delta^{AE} \delta^{b'c'} \delta^{d'f'} + \delta^{AE} \delta^{bf} \left(5 \delta^{cd} + 3 \delta^{c'd'} \right) \\
& \left. - 2 \delta^{CD} \left[9 \left(\delta^{BE} \delta^{AF} + \delta^{AE} \delta^{BF} \right) - \delta^{be} \delta^{af} \right. \right. \\
& \left. \left. - \delta^{ae} \delta^{bf} + \delta^{ab} \delta^{ef} + \delta^{b'e'} \delta^{a'f'} + \delta^{a'e'} \delta^{b'f'} - \delta^{a'b'} \delta^{e'f'} \right] \right\}. \quad (4.1.19)
\end{aligned}$$

This matrix element exhibits the same antisymmetry between the $SO(4)_{\text{AdS}}$ and $SO(4)_{S^5}$ indices that is exhibited in eqn. (4.1.3). The off-diagonal H_{BF} mixing sector is essentially equivalent to its counterpart in eqn. (4.1.4):

$$\langle J | a_n^A a_n^B a_{-2n}^C (H_{\text{BF}}) a_{-2n}^{D\dagger} b_n^{\alpha\dagger} b_n^{\beta\dagger} | J \rangle = \frac{n^2 \lambda'}{2J} \delta^{CD} \left\{ \left(\gamma^{ab'} \right)^{\alpha\beta} - \left(\gamma^{a'b} \right)^{\alpha\beta} \right\}. \quad (4.1.20)$$

The diagonal contributions from the pure fermion sector H_{FF} in the (2, 2) and (3, 3) blocks of table 4.11 appear as

$$\begin{aligned} \left\langle J \left| b_n^\alpha b_n^\beta a_{-2n}^A (H_{\text{FF}}) a_{-2n}^{B\dagger} b_n^{\gamma\dagger} b_n^{\delta\dagger} \right| J \right\rangle &= \frac{n^2 \lambda'}{24J} \delta^{AB} \left\{ (\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\beta\delta} - (\gamma^{ij})^{\alpha\beta} (\gamma^{ij})^{\gamma\delta} \right. \\ &\quad \left. - (\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\beta\gamma} - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\beta\delta} + (\gamma^{i'j'})^{\alpha\beta} (\gamma^{i'j'})^{\gamma\delta} + (\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\beta\gamma} \right\}. \end{aligned} \quad (4.1.21)$$

The H_{BF} sector exhibits the following contribution to the lower diagonal blocks (2, 2) and (3, 3):

$$\begin{aligned} \left\langle J \left| b_n^\alpha b_n^\beta a_{-2n}^A (H_{\text{BF}}) a_{-2n}^{B\dagger} b_n^{\gamma\dagger} b_n^{\delta\dagger} \right| J \right\rangle &= \frac{n^2 \lambda'}{J} \left\{ -10 \delta^{a'b'} (\delta^{\alpha\delta} \delta^{\beta\gamma} - \delta^{\alpha\gamma} \delta^{\beta\delta}) \right. \\ &\quad - 8 \delta^{ab} (\delta^{\alpha\delta} \delta^{\beta\gamma} - \delta^{\alpha\gamma} \delta^{\beta\delta}) - \delta^{\alpha\gamma} [(\gamma^{ab})^{\beta\delta} - (\gamma^{a'b'})^{\beta\delta}] + \delta^{\alpha\delta} [(\gamma^{ab})^{\beta\gamma} - (\gamma^{a'b'})^{\beta\gamma}] \\ &\quad \left. + \delta^{\beta\gamma} [(\gamma^{ab})^{\alpha\delta} - (\gamma^{a'b'})^{\alpha\delta}] - \delta^{\beta\delta} [(\gamma^{ab})^{\alpha\gamma} - (\gamma^{a'b'})^{\alpha\gamma}] \right\}. \end{aligned} \quad (4.1.22)$$

Finally, the off-diagonal version of (4.1.22) appears in the (2, 3) block (along with its transpose in the (3, 2) block):

$$\begin{aligned} \left\langle J \left| b_n^\alpha b_n^\beta a_{-2n}^A (H_{\text{BF}}) a_n^{B\dagger} b_n^{\gamma\dagger} b_{-2n}^{\delta\dagger} \right| J \right\rangle &= -\frac{n^2 \lambda'}{J} \left\{ \delta^{a'b'} (\delta^{\alpha\delta} \delta^{\beta\gamma} - \delta^{\alpha\gamma} \delta^{\beta\delta}) \right. \\ &\quad \left. - \delta^{ab} (\delta^{\alpha\delta} \delta^{\beta\gamma} - \delta^{\alpha\gamma} \delta^{\beta\delta}) + \delta^{\alpha\gamma} [(\gamma^{ab})^{\beta\delta} - (\gamma^{a'b'})^{\beta\delta}] - \delta^{\beta\gamma} [(\gamma^{ab})^{\alpha\delta} - (\gamma^{a'b'})^{\alpha\delta}] \right\}. \end{aligned} \quad (4.1.23)$$

The fermionic sector perturbation matrix is displayed schematically in table 4.12. Like table 4.11, it is $1,024 \times 1,024$ once redundant index assignments are eliminated.

H_{int}	$b_{-2n}^{\zeta\dagger} b_n^{\epsilon\dagger} b_n^{\delta\dagger} J\rangle$	$a_{-2n}^{C\dagger} a_n^{D\dagger} b_n^{\delta\dagger} J\rangle$	$a_n^{C\dagger} a_n^{D\dagger} b_{-2n}^{\delta\dagger} J\rangle$
$\langle J b_n^\alpha b_n^\beta b_{-2n}^\gamma$	H_{FF}	H_{BF}	H_{BF}
$\langle J b_n^\alpha a_n^A a_{-2n}^B$	H_{BF}	$H_{\text{BB}} + H_{\text{BF}}$	H_{BF}
$\langle J b_{-2n}^\alpha a_n^A a_n^B$	H_{BF}	H_{BF}	$H_{\text{BB}} + H_{\text{BF}}$

Table 4.12: Fermionic string perturbation matrix ($q = r = n$, $s = -2n$)

The purely fermionic subsector in the $(1, 1)$ block of table 4.12 takes the form

$$\begin{aligned}
\left\langle J \left| b_n^\alpha b_n^\beta b_{-2n}^\gamma (H_{\text{FF}}) b_{-2n}^{\zeta\dagger} b_n^{\epsilon\dagger} b_n^{\delta\dagger} \right| J \right\rangle &= \frac{9n^2 \lambda'}{J} \delta^{\gamma\zeta} (\delta^{\alpha\epsilon} \delta^{\beta\delta} - \delta^{\alpha\delta} \delta^{\beta\epsilon}) \\
&+ \frac{n^2 \lambda'}{24J} \left\{ \delta^{\gamma\zeta} \left[(\gamma^{ij})^{\alpha\beta} (\gamma^{ij})^{\delta\epsilon} - (\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\beta\epsilon} + (\gamma^{ij})^{\alpha\epsilon} (\gamma^{ij})^{\beta\delta} - (\gamma^{i'j'})^{\alpha\beta} (\gamma^{i'j'})^{\delta\epsilon} \right. \right. \\
&+ (\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\beta\epsilon} - (\gamma^{i'j'})^{\alpha\epsilon} (\gamma^{i'j'})^{\beta\delta} \left. \right] - 2\delta^{\alpha\delta} \left[(\gamma^{ij})^{\beta\gamma} (\gamma^{ij})^{\epsilon\zeta} - (\gamma^{ij})^{\beta\epsilon} (\gamma^{ij})^{\gamma\zeta} \right. \\
&+ (\gamma^{ij})^{\beta\zeta} (\gamma^{ij})^{\gamma\epsilon} - (\gamma^{i'j'})^{\beta\gamma} (\gamma^{i'j'})^{\epsilon\zeta} + (\gamma^{i'j'})^{\beta\epsilon} (\gamma^{i'j'})^{\gamma\zeta} - (\gamma^{i'j'})^{\beta\zeta} (\gamma^{i'j'})^{\gamma\epsilon} \left. \right] \\
&+ 2\delta^{\alpha\epsilon} \left[(\gamma^{ij})^{\beta\gamma} (\gamma^{ij})^{\delta\zeta} - (\gamma^{ij})^{\beta\delta} (\gamma^{ij})^{\gamma\zeta} + (\gamma^{ij})^{\beta\zeta} (\gamma^{ij})^{\gamma\delta} - (\gamma^{i'j'})^{\beta\gamma} (\gamma^{i'j'})^{\delta\zeta} \right. \\
&+ (\gamma^{i'j'})^{\beta\delta} (\gamma^{i'j'})^{\gamma\zeta} - (\gamma^{i'j'})^{\beta\zeta} (\gamma^{i'j'})^{\gamma\delta} \left. \right] + 2\delta^{\beta\delta} \left[(\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\epsilon\zeta} - (\gamma^{ij})^{\alpha\epsilon} (\gamma^{ij})^{\gamma\zeta} \right. \\
&+ (\gamma^{ij})^{\alpha\zeta} (\gamma^{ij})^{\gamma\epsilon} - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\epsilon\zeta} + (\gamma^{i'j'})^{\alpha\epsilon} (\gamma^{i'j'})^{\gamma\zeta} - (\gamma^{i'j'})^{\alpha\zeta} (\gamma^{i'j'})^{\gamma\epsilon} \left. \right] \\
&- 2\delta^{\beta\epsilon} \left[(\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\delta\zeta} - (\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\gamma\zeta} + (\gamma^{ij})^{\alpha\zeta} (\gamma^{ij})^{\gamma\delta} - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\delta\zeta} \right. \\
&+ (\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\gamma\zeta} - (\gamma^{i'j'})^{\alpha\zeta} (\gamma^{i'j'})^{\gamma\delta} \left. \right] \left. \right\}. \tag{4.1.24}
\end{aligned}$$

The off-diagonal blocks $(1, 2)$ and $(1, 3)$ receive contributions from the H_{BF} sector:

$$\begin{aligned}
\left\langle J \left| b_n^\alpha b_n^\beta b_{-2n}^\gamma (H_{\text{BF}}) a_{-2n}^{A\dagger} a_n^{B\dagger} b_n^{\delta\dagger} \right| J \right\rangle &= \frac{n^2 \lambda'}{J} \left\{ \delta^{\alpha\delta} \left[(\gamma^{ab'})^{\beta\gamma} - (\gamma^{a'b})^{\beta\gamma} \right] \right. \\
&\left. - \delta^{\delta\beta} \left[(\gamma^{ab'})^{\alpha\gamma} - (\gamma^{a'b})^{\alpha\gamma} \right] \right\}. \tag{4.1.25}
\end{aligned}$$

The bosonic sector H_{BB} contributes to the (2, 2) and (3, 3) blocks:

$$\begin{aligned} \langle J | b_q^\alpha a_r^A a_s^B (H_{\text{BB}}) a_s^{C\dagger} a_r^{D\dagger} b_q^{\beta\dagger} | J \rangle &= -\frac{n^2 \lambda'}{2J} \delta^{\alpha\beta} \left\{ 9 \delta^{AD} \delta^{BC} + 4 \delta^{ac} \delta^{bd} - 4 \delta^{ab} \delta^{cd} \right. \\ &\quad \left. - \delta^{ad} (5 \delta^{bc} + 3 \delta^{b'c'}) - 4 \delta^{a'c'} \delta^{b'd'} + 4 \delta^{a'b'} \delta^{c'd'} + \delta^{a'd'} (5 \delta^{b'c'} + 3 \delta^{bc}) \right\}. \end{aligned} \quad (4.1.26)$$

In the same lower-diagonal blocks, H_{BF} exhibits the contribution

$$\begin{aligned} \langle J | b_n^\alpha a_n^A a_{-2n}^B (H_{\text{BF}}) a_{-2n}^{C\dagger} a_n^{D\dagger} b_n^{\beta\dagger} | J \rangle &= -\frac{n^2 \lambda'}{8J} \left\{ 39 \delta^{\alpha\beta} \delta^{AD} \delta^{BC} \right. \\ &\quad + \delta^{\alpha\beta} \delta^{AD} (\delta^{b'c'} - 7 \delta^{bc}) - 4 \delta^{\alpha\beta} \delta^{BC} (\delta^{ad} - \delta^{a'd'}) + 4 \delta^{BC} [(\gamma^{ad})^{\alpha\beta} - (\gamma^{a'd'})^{\alpha\beta}] \\ &\quad \left. - 8 \delta^{AD} [(\gamma^{bc})^{\alpha\beta} - (\gamma^{b'c'})^{\alpha\beta}] \right\}. \end{aligned} \quad (4.1.27)$$

Finally, H_{BF} yields matrix elements in the off-diagonal block (2, 3):

$$\begin{aligned} \langle J | b_n^\alpha a_n^A a_{-2n}^B (H_{\text{BF}}) a_n^{C\dagger} a_n^{D\dagger} b_{-2n}^{\beta\dagger} | J \rangle &= -\frac{n^2 \lambda'}{32J} \left\{ 9 \delta^{\alpha\beta} \delta^{AC} \delta^{BD} + 9 \delta^{\alpha\beta} \delta^{AD} \delta^{BC} \right. \\ &\quad + \delta^{\alpha\beta} \delta^{AC} (23 \delta^{bd} - 41 \delta^{b'd'}) + \delta^{\alpha\beta} \delta^{AD} (23 \delta^{bc} - 41 \delta^{b'c'}) \\ &\quad \left. - 32 \delta^{AD} [(\gamma^{bc})^{\alpha\beta} - (\gamma^{b'c'})^{\alpha\beta}] - 32 \delta^{AC} [(\gamma^{bd})^{\alpha\beta} - (\gamma^{b'd'})^{\alpha\beta}] \right\}. \end{aligned} \quad (4.1.28)$$

We can perform a full symbolic diagonalization of the $1,024 \times 1,024$ bosonic and fermionic perturbation matrices to obtain the one-loop in λ' , $O(1/J)$ energy corrections. They can all be expressed in terms of dimensionless eigenvalues Λ according to the standard formula (4.1.13) modified by setting $q = r = n$, $s = -2n$:

$$E_J(n) = 3 + 3n^2 \lambda' \left(1 + \frac{\Lambda}{J} + O(J^{-2}) \right). \quad (4.1.29)$$

The resulting spectrum is displayed in table 4.13. The levels clearly organize themselves into two superconformal multiplets built on vector primary states. Note that the spinor multiplet is absent and that the degeneracy between multiplets that was

Λ_1 (S^5 vector)	$-23/3$	$-20/3$	$-17/3$	$-14/3$	$-11/3$	$-8/3$	$-5/3$	$-2/3$	$1/3$
Multiplicity	4_B	32_F	112_B	224_F	280_B	224_F	112_B	32_F	4_B
Λ_2 (AdS_5 vector)	$-19/3$	$-16/3$	$-13/3$	$-10/3$	$-7/3$	$-4/3$	$-1/3$	$2/3$	$5/3$
Multiplicity	4_B	32_F	112_B	224_F	280_B	224_F	112_B	32_F	4_B

Table 4.13: Spectrum of three-impurity string Hamiltonian with $(q = r = n, s = -2n)$

seen in the inequivalent mode index case has been lifted. The spinor multiplet is absent for the following reason: it contains a representation at level $L = 4$ arising from fermion creation operators completely symmetrized on $SO(4) \times SO(4)$ spinor indices; such a construct must vanish unless all the creation operator mode indices are different.

If we keep track of the $SO(4) \times SO(4)$ irrep structure, we find that the symmetric-traceless bosonic $SO(4)_{S^5}$ states arising from the closed $\mathfrak{su}(2)$ subsector fall into the $-11/3$ [280_B] level. This is the counterpart of the -4 [280_B] level in table 4.10. To compare with Minahan and Zarembo's Bethe ansatz calculation of the corresponding gauge theory operator dimension, we must evaluate eqn. (4.1.17) with the appropriate choice of parameters. In particular, $M_n = 2$ when two mode indices are allowed to coincide and, comparing with eqn. (4.1.29), we find perfect agreement with the string theory prediction $\Lambda = -11/3$. States at level $L = 4$ in the second multiplet in table 4.13 correspond to operators in the $\mathfrak{sl}(2)$ closed sector of the gauge theory and the eigenvalue $\Lambda = -7/3$ [280_B] amounts to a prediction for the one-loop anomalous dimension of that class of gauge theory operators. As mentioned at the end of the previous subsection, we will need to develop a numerical treatment of the $\mathfrak{sl}(2)$ spin chain Hamiltonian in order to assess the agreement between the string theory and gauge theory in this sector.

4.2 Three-impurity spectrum: all orders in λ'

In the previous section, we have studied the eigenvalue spectrum of the string theory perturbation Hamiltonian expanded to leading order in $1/J$ and to one-loop order in λ' . The expansion in λ' was for convenience only since our expressions for matrix elements are exact in this parameter. We should, in principle, be able to obtain results that are exact in λ' (but still of leading order in $1/J$). This is a worthwhile enterprise since recent progress on the gauge theory side has made it possible to evaluate selected operator anomalous dimensions to two- and three-loop order. The simple one-loop calculations of the previous sections have given us an overview of how the perturbed string theory eigenvalues are organized into superconformal multiplets. This provides a very useful orientation for the more complex all-orders calculation, to which we now turn.

4.2.1 Inequivalent mode indices: $(q \neq r \neq s)$

Our first step is to collect the exact matrix elements of the perturbing Hamiltonian between three-impurity states of unequal mode indices. The block structure of the perturbation matrix in the spacetime boson sector is given in table 4.2 and the exact form of the $(1, 1)$ block is

$$\begin{aligned} \langle J | a_q^A a_r^B a_s^C (H_{\text{BB}}) a_s^{D\dagger} a_r^{E\dagger} a_q^{F\dagger} | J \rangle = & -\frac{1}{2\omega_q \omega_r \omega_s} \left\{ \delta^{BE} \omega_r \left[\delta^{CD} \delta^{AF} (s^2 + q^2(1 + 2s^2 \lambda')) \right. \right. \\ & - (q^2 + s^2) \delta^{cd} \delta^{af} - 2qs(\delta^{ad} \delta^{cf} - \delta^{ac} \delta^{df}) + (q^2 - s^2) \delta^{af} \delta^{c'd'} - (q^2 - s^2) \delta^{a'f'} \delta^{cd} \\ & + (q^2 + s^2) \delta^{c'd'} \delta^{a'f'} + 2qs(\delta^{a'd'} \delta^{c'f'} - \delta^{a'c'} \delta^{d'f'}) \left. \right] + (C \rightleftharpoons B, D \rightleftharpoons E, s \rightleftharpoons r) \\ & \left. + (A \rightleftharpoons B, F \rightleftharpoons E, q \rightleftharpoons r) \right\}, \end{aligned} \quad (4.2.1)$$

where we define $\omega_q \equiv \sqrt{q^2 + 1/\lambda'}$ to simplify this and other similar expressions.

The off-diagonal H_{BF} contributions to the $(1, 2)$, $(1, 3)$ and $(1, 4)$ blocks are yet

more complicated. To simplify the expressions, we define

$$\begin{aligned} F_1 &\equiv \sqrt{(\omega_q + q)(\omega_r - r)} , & F_2 &\equiv \sqrt{(\omega_q - q)(\omega_r + r)} , \\ F_3 &\equiv \sqrt{(\omega_q - q)(\omega_r - r)} , & F_4 &\equiv \sqrt{(\omega_q + q)(\omega_r + r)} . \end{aligned} \quad (4.2.2)$$

Using these functions, the matrix elements in these off-diagonal subsectors are given by:

$$\begin{aligned} \langle J | a_q^A a_r^B a_s^C (H_{\text{BF}}) a_s^{D\dagger} b_r^{\alpha\dagger} b_q^{\beta\dagger} | J \rangle = & \frac{\delta^{CD}}{32\omega_q\omega_r J} \left\{ \frac{8}{\sqrt{\lambda'}} (F_1 - F_2) \delta^{AB} \delta^{\gamma\delta} \right. \\ & - 2(q - r)(F_3 + F_4) \delta^{AB} \delta^{\gamma\delta} + 4(q - r)(F_3 + F_4) (\gamma^{ab})^{\gamma\delta} \\ & - 2(q + r)(F_3 - F_4) (\gamma^{ab'})^{\gamma\delta} + (2qF_3 - 2qF_4 + 2rF_3 - 2rF_4) (\gamma^{a'b})^{\gamma\delta} \\ & - (4qF_3 + 4qF_4 - 4rF_3 - 4rF_4) (\gamma^{a'b'})^{\gamma\delta} + \frac{8}{\sqrt{\lambda'}} (F_2 - F_1) \delta^{a'b'} \delta^{\gamma\delta} \\ & + 4(q - r)(F_3 + F_4) \delta^{\gamma\delta} \delta^{a'b'} - 2(q - r)(F_3 + F_4) \delta^{\gamma\delta} (\delta^{ab} - \delta^{a'b'}) \\ & - 4\lambda' \omega_q \omega_r (q - r)(F_3 + F_4) \delta^{AB} \delta^{\gamma\delta} \\ & + 4\sqrt{\lambda'} (qr - \omega_q \omega_r) \left[(F_1 + F_2) \left((\gamma^{ab'})^{\gamma\delta} - (\gamma^{a'b})^{\gamma\delta} \right) - (F_1 - F_2) \delta^{\gamma\delta} (\delta^{ab} - \delta^{a'b'}) \right] \\ & + 2(\omega_q + \omega_r)(F_3 + F_4) \left[(\gamma^{ab'})^{\gamma\delta} - (\gamma^{a'b})^{\gamma\delta} \right] \\ & + 4\sqrt{\lambda'} (r\omega_q - q\omega_r)(F_1 + F_2) \left[(\gamma^{ab})^{\gamma\delta} - (\gamma^{a'b'})^{\gamma\delta} \right] \\ & - 4\lambda' (q - r)(F_3 - F_4)(r\omega_q + q\omega_r) \delta^{AB} \delta^{\gamma\delta} - \lambda' \delta^{AB} \\ & + 2\lambda' (\omega_q \omega_r - qr)(q - r)(F_3 + F_4) \delta^{AB} \delta^{\gamma\delta} + 4\sqrt{\lambda'} (\omega_q \omega_r + qr)(F_1 - F_2) \delta^{AB} \delta^{\gamma\delta} \\ & \left. - 2\lambda' (q - r)(\omega_q \omega_r + qr)(F_3 + F_4) \delta^{AB} \delta^{\gamma\delta} \right\} . \end{aligned} \quad (4.2.3)$$

The H_{FF} contribution to the lower-diagonal blocks (2, 2), (3, 3) and (4, 4) is

$$\begin{aligned} \langle J | b_q^{\alpha} b_r^{\beta} a_s^A (H_{\text{FF}}) a_s^{B\dagger} b_r^{\gamma\dagger} b_q^{\delta\dagger} | J \rangle = & \frac{\delta^{AB}}{48\omega_r\omega_s J} \sqrt{\lambda'} \left\{ 2rs \sqrt{1/\lambda'} \left[\left((\gamma^{ij})^{\alpha\gamma} (\gamma^{ij})^{\beta\delta} - (\gamma^{i'j'})^{\alpha\gamma} (\gamma^{i'j'})^{\beta\delta} \right) \right. \right. \\ & \left. \left. - \left((\gamma^{ij})^{\alpha\delta} (\gamma^{ij})^{\beta\gamma} - (\gamma^{i'j'})^{\alpha\delta} (\gamma^{i'j'})^{\beta\gamma} \right) - \left((\gamma^{ij})^{\alpha\beta} (\gamma^{ij})^{\gamma\delta} - (\gamma^{i'j'})^{\alpha\beta} (\gamma^{i'j'})^{\gamma\delta} \right) \right] \right\} \end{aligned}$$

$$-12 \left[2\delta^{\alpha\delta}\delta^{\beta\gamma} \left(s^2\sqrt{1/\lambda'} - 2rs\sqrt{\lambda'}\omega_r\omega_s + r^2(2s^2\sqrt{\lambda'} + \sqrt{1/\lambda'}) \right) \right] \Big\} . \quad (4.2.4)$$

The bose-fermi Hamiltonian H_{BF} contributes the following matrix elements to the same lower-diagonal blocks:

$$\begin{aligned} \langle J | b_q^\alpha b_r^\beta a_s^A (H_{\text{BF}}) a_s^{B\dagger} b_r^{\gamma\dagger} b_q^{\delta\dagger} | J \rangle = & -\frac{1}{2\omega_q\omega_r\omega_s} \left\{ s\sqrt{\lambda'}\delta^{ab}\delta^{\alpha\delta}\delta^{\beta\gamma} \left[s\omega_r(2q^2\sqrt{\lambda'} + \sqrt{1/\lambda'}) \right. \right. \\ & + s\omega_q(2r^2\sqrt{\lambda'} + \sqrt{1/\lambda'}) - 2\omega_q\omega_r\omega_s(q+r)\sqrt{\lambda'} \Big] + \delta^{a'b'}\delta^{\alpha\delta}\delta^{\beta\gamma} \left[2\omega_rq^2(1+s^2\lambda') \right. \\ & + s^2\omega_r + 2\omega_qr^2(1+s^2\lambda') + s^2\omega_q - 2s(q+r)\lambda'\omega_q\omega_r\omega_s \Big] \\ & \left. + sr\omega_q\delta^{\alpha\delta} \left[(\gamma^{ab})^{\beta\gamma} - (\gamma^{a'b'})^{\beta\gamma} \right] + sq\omega_r\delta^{\beta\gamma} \left[(\gamma^{ab})^{\alpha\delta} - (\gamma^{a'b'})^{\alpha\delta} \right] \right\} . \quad (4.2.5) \end{aligned}$$

To simplify off-diagonal elements in the (2, 3), (2, 4) and (3, 4) blocks, we define

$$\begin{aligned} G_1 &\equiv \sqrt{(\omega_r + r)(\omega_s - s)} , & G_2 &\equiv \sqrt{(\omega_r - r)(\omega_s + s)} , \\ G_3 &\equiv \sqrt{(\omega_r - r)(\omega_s - s)} , & G_4 &\equiv \sqrt{(\omega_r + r)(\omega_s + s)} . \end{aligned} \quad (4.2.6)$$

The matrix elements in these subsectors are then given by

$$\begin{aligned} \langle J | b_q^\alpha b_r^\beta a_s^A (H_{\text{BF}}) a_r^{B\dagger} b_q^{\gamma\dagger} b_s^{\delta\dagger} | J \rangle = & -\frac{1}{16(\lambda'\omega_r\omega_s)^{3/2}} \left\{ \sqrt{\omega_r\omega_s}\lambda'\delta^{\alpha\gamma} \left[2\delta^{ab}\delta^{\beta\delta} \left[(G_1 + G_2)(2 - 2\lambda'\omega_r\omega_s) \right. \right. \right. \\ & + (r+s)\sqrt{\lambda'} \left(G_4 - \lambda'G_4(r - \omega_r)(s - \omega_s) \right. \\ & \left. \left. + G_3(-1 + rs\lambda' + r\omega_s\lambda' + \omega_r(s + \omega_s)\lambda') \right) \right] \\ & + 2\sqrt{\lambda'} \left[(r+s)(G_3 - G_4) + \sqrt{\lambda'}(G_1 - G_2)(r\omega_s - s\omega_r) \right] \left[(\gamma^{ab})^{\beta\delta} - (\gamma^{a'b'})^{\beta\delta} \right] \\ & + \sqrt{\lambda'} \left[2rs\sqrt{\lambda'}G_1 - 2rs\sqrt{\lambda'}G_2 + (r-s)(G_3 + G_4) + (\omega_s - \omega_r)(G_3 - G_4) \right. \\ & \left. + 2\omega_r\omega_s\sqrt{\lambda'}(G_2 - G_1) \right] \left[(\gamma^{ab'})^{\beta\delta} - (\gamma^{a'b})^{\beta\delta} \right] + 2\delta^{a'b'}\delta^{\beta\delta} \left[-2rs\sqrt{\lambda'}(G_1 - G_2) \right. \\ & \left. + (r+s)\sqrt{\lambda'} \left(-G_4 - \lambda'G_4(r - \omega_r)(s - \omega_s) \right. \right. \\ & \left. \left. + G_3(1 + rs\lambda' + r\omega_s\lambda' + \omega_r(s + \omega_s)\lambda') \right) \right] \Big\} . \quad (4.2.7) \end{aligned}$$

The entries in the spacetime fermion block matrix of table 4.7 are far too complicated to write out explicitly: they are best generated, viewed and manipulated with computer algebra techniques. The explicit formulas, along with a collection of the Mathematica programs written to generate and work with them, are available on the web.³

We were not able to symbolically diagonalize the complete perturbation matrix built from the exact (in λ') matrix elements listed above: with the computing resources available to us, the routines for diagonalizing the full 2,048-dimensional matrices would not terminate in any reasonable time. As noted in the previous section, however, gauge theory arguments suggest that there are three protected $SO(4) \times SO(4)$ irreps that do not mix with any other irreps. It is a straightforward matter to project the perturbation matrix onto these unique protected irreps to obtain analytic expressions for the corresponding exact eigenvalues. In fact, the superconformal multiplet structure of the three-impurity problem is such that the energies/dimensions of all other irreps can be inferred from those of the three protected irreps. Hence, this method will give us exact expressions for all the energy levels of the three-impurity problem.

Consider first the $\mathfrak{sl}(2)$ closed sector. The dual sector is generated on the string theory side by bosonic creation operators completely symmetrized (and traceless) on $SO(4)_{AdS}$ vector indices. The simplest way to make this projection on eqn. (4.2.1) is to compute diagonal elements between the symmetrized states

$$a_q^{(a\dagger} a_r^{b\dagger} a_s^{c\dagger)} |J\rangle \ , \quad (4.2.8)$$

with $a \neq b \neq c$ (and, of course, $a, b, c \in 1, \dots, 4$). The charges of the fermionic oscillators under this subgroup are $\pm 1/2$, so the three-boson state of this type cannot mix with one boson and two fermions (or any other state). Hence, the above projection

³<http://theory.caltech.edu/~swanson/MMA1/mma1.html>

of eqn. (4.2.1) yields the closed sector eigenvalue correction

$$\begin{aligned} \delta E_{AdS}(q, r, s, J) &= \frac{1}{J\omega_q\omega_r\omega_s} \left\{ qs(1 - qs\lambda')\omega_r + qr(1 - qr\lambda')\omega_s + rs(1 - rs\lambda')\omega_q \right. \\ &\quad \left. + [qr + s(q + r)]\lambda'\omega_q\omega_r\omega_s \right\} \\ &\approx \frac{1}{J} \left\{ -2(q^2 + qr + r^2)\lambda' - \frac{15}{8}(q^2r^2(q + r)^2)\lambda'^3 + \dots \right\}. \end{aligned} \quad (4.2.9)$$

To facilitate eventual comparison with gauge theory results, we have performed a small- λ' expansion in the final line with the substitution $s \rightarrow -(q + r)$ (since the mode indices satisfy the constraint $s + q + r = 0$). The leading correction $-2(q^2 + qr + r^2)\lambda'$ reproduces the one-loop eigenvalue $\Lambda_{BB} = -2$ [280_B] located at level $L = 4$ in the $SO(4)_{AdS}$ multiplet in table 4.10.

The closed $\mathfrak{su}(2)$ sector is generated by bosonic creation operators completely symmetrized on traceless $SO(4)_{S^5}$ indices. Projection onto this irrep is most simply achieved by choosing all mode operators in eqn. (4.2.1) to carry symmetrized, traceless $SO(4)_{S^5}$ labels (they can also be thought of as carrying charge +1 under some $SO(2)$ subgroup of $SO(4)_{S^5}$). Direct projection yields the $SO(4)_{S^5}$ eigenvalue

$$\begin{aligned} \delta E_{S^5}(q, r, s, J) &= -\frac{1}{J\omega_q\omega_r\omega_s} \left\{ [qr + r^2 + q^2(1 + r^2\lambda')]\omega_s \right. \\ &\quad + [qs + s^2 + q^2(1 + s^2\lambda')]\omega_r \\ &\quad \left. + [rs + s^2 + r^2(1 + s^2\lambda')]\omega_q - [rs + q(r + s)]\lambda'\omega_q\omega_r\omega_s \right\} \\ &\approx \frac{1}{J} \left\{ -4(q^2 + qr + r^2)\lambda' + (q^2 + qr + r^2)^2\lambda'^2 \right. \\ &\quad \left. - \frac{3}{4}(q^6 + 3q^5r + 8q^4r^2 + 11q^3r^3 + 8q^2r^4 + 3qr^5 + r^6)\lambda'^3 + \dots \right\}. \end{aligned} \quad (4.2.10)$$

This is the all-loop formula corresponding to gauge theory operator dimensions in the closed $\mathfrak{su}(2)$ subsector; the leading-order term $-4(q^2 + qr + r^2)\lambda'$ reproduces the one-loop eigenvalue $\Lambda_{BB} = -4$ [280_B] at level $L = 4$ in the $SO(4)_{S^5}$ vector multiplet in table 4.10.

The eigenvalue of the symmetrized pure-fermion irrep can be obtained by evaluating the exact matrix element H_{FF} acting on three symmetrized fermionic creation operators with $SO(4) \times SO(4)$ indices chosen to lie in the same Π projection (with inequivalent mode indices). The exact energy shift for this irrep turns out to be

$$\begin{aligned} \delta E_{\text{Fermi}}(q, r, s, J) &= -\frac{1}{4J\omega_q\omega_r\omega_s} \left\{ -4(rs + q(r+s))\lambda'\omega_q\omega_r\omega_s \right. \\ &\quad \left. + \left[\omega_q(2s^2 + 4r^2s^2\lambda' + 2r^2) + (s \rightarrow r, r \rightarrow q, q \rightarrow s) + (q \Rightarrow r) \right] \right\} \\ &\approx \frac{1}{J} \left\{ -3(q^2 + qr + r^2)\lambda' + \frac{1}{2}(q^2 + qr + r^2)^2\lambda'^2 \right. \\ &\quad \left. - \frac{3}{16}(2q^6 + 6q^5r + 21q^4r^2 + 32q^3r^3 + 21q^2r^4 + 6qr^5 + 2r^6)\lambda'^3 + \dots \right\}. \quad (4.2.11) \end{aligned}$$

The leading-order λ' correction $-3(q^2 + qr + r^2)\lambda'$ reproduces the $\Lambda_{\text{FF}} = -3$ [580_F] eigenvalue at the $L = 4$ level in the spinor multiplet in table 4.10. This and the higher order terms in the eigenvalue will eventually be compared with the dimensions of operators in the closed, fermionic $\mathfrak{su}(1|1)$ sector in the gauge theory.

The argument we are making relies heavily on the claim that the perturbation matrix is block diagonal on the closed subsectors described above: we have evaluated the exact energy shift on these subsectors by simply taking the diagonal matrix element of the perturbing Hamiltonian in a particular state in each sector. We will now carry out a simple numerical test of the claimed block diagonalization of the full perturbing Hamiltonian. The basic idea is that, while it is impractical to algebraically diagonalize the full $2,048 \times 2,048$ perturbation matrices, it is quite easy to do a numerical diagonalization for a specific choice of λ' and mode indices q, r, s . One can then check that the numerical eigenvalues match the analytic predictions evaluated at the chosen coupling and mode indices. For definiteness, we choose

$$q = 1, \quad r = 2, \quad s = -3, \quad \lambda' = 1. \quad (4.2.12)$$

The predicted eigenvalue shifts of the three protected states, evaluated at the pa-

parameter choices of (4.2.12) are given in table 4.14. These values come directly from eqns. (4.2.9, 4.2.10, 4.2.11) above (with J set to unity, for convenience). Since we

$\delta E : \lambda' = 1$	$q = 1, r = 2, s = -3$
$\delta E_{AdS}(1, 2, -3, J = 1)$	$= -16.255434067000426$
$\delta E_{S^5}(1, 2, -3, J = 1)$	$= -20.137332508389193$
$\delta E_{\text{Fermi}}(1, 2, -3, J = 1)$	$= -18.19638328769481$

Table 4.14: Exact numerical eigenvalues of three-impurity protected sectors

want to compare these energies to a numerical diagonalization, we must maintain a high level of precision in the numerical computation. With the parameter choices of (4.2.12), the numerical diagonalization of the full $2,048 \times 2,048$ perturbation matrices on both the spacetime boson (table 4.2) and spacetime fermion (table 4.3) sectors yields the spectrum and multiplicities displayed in table 4.15. The multiplicities are consistent with the superconformal multiplet structure we found in the one-loop analysis (given in table 4.10). The predicted closed sector eigenvalues (listed in table 4.14) match, to the precision of the calculation, entries in the list of numerical eigenvalues. These energies also appear at the expected levels within the multiplets. $E_{AdS}(1, 2, -3, J)$ and $E_{S^5}(1, 2, -3, J)$ appear in bosonic levels with multiplicity 280_B , while energy $E_{\text{Fermi}}(1, 2, -3, J)$ appears as a fermionic level with multiplicity 560_F ; according to table 4.10 these are uniquely identified as the central $L = 4$ levels of their respective multiplets, exactly where the protected energy levels must lie. All of this is clear evidence that the closed sector states of the string theory do not mix with other states under the perturbing Hamiltonian, thus justifying our method of calculating their exact eigenenergies.

At one loop, we found that the three superconformal multiplets were displaced from each other by precisely the internal level spacing. This led to an accidental degeneracy that is lifted in the exact dimension formulas we have just derived. To explore this, it is useful to have formulas for the eigenvalues of all the levels in each multiplet. From the discussion in Section 4.1, we see that each level in the string

$\delta E(1, 2, -3, J = 1) \lambda' = 1$	Mult.	$\delta E(1, 2, -3, J = 1) \lambda' = 1$	Mult.
-30.821354623065	4_B	-28.8804054023706	8_F
-26.9394561816763	4_B	-28.150349094396	32_F
-26.2093998737015	64_B	-24.2684506530072	32_F
-25.4793435657269	112_B	-23.5383943450326	224_F
-21.5974451243382	112_B	-22.808338037058	224_F
-20.8673888163637	448_B	-18.9264395956693	224_F
-20.1373325083891	280_B	-18.1963832876947	560_F
-16.2554340670003	280_B	-17.4663269797201	224_F
-15.5253777590258	448_B	-13.5844285383314	224_F
-14.7953214510512	112_B	-12.8543722303568	224_F
-10.9134230096624	112_B	-12.1243159223822	32_F
-10.1833667016878	64_B	-8.24241748099347	32_F
-9.4533103937133	4_B	-7.51236117301893	8_F
-5.57141195232456	4_B		

Table 4.15: All-loop numerical spectrum of three-impurity states ($q = 1$, $r = 2$, $s = -3$, $\lambda' = 1$, $J = 1$). Left panel: bosons; right panel: fermions

energy spectrum can be connected by a simple integer shift in the angular momentum J . Since we are working at $O(1/J)$ in a large- J expansion, all contributions from this shift must come from the BMN limit of the theory. In other words, by sending $J \rightarrow J + 2 - L/2$ in the BMN formula for the energy

$$E = \sqrt{1 + \frac{n^2 g_{YM}^2 N_c}{(J + 2 - L/2)^2}} + \dots, \quad (4.2.13)$$

we can generate an expansion, to arbitrary order in λ' , for each level L in the entire superconformal multiplet.

For the vector $SO(4)_{AdS}$ multiplet, we find

$$\begin{aligned}
\delta E_{AdS}(q, r, J, L) \approx & \frac{1}{J} \left\{ (L-6)(q^2 + qr + r^2)\lambda' - \frac{1}{2}(L-4)(q^2 + qr + r^2)^2\lambda'^2 \right. \\
& + \frac{3}{16} \left[2(L-4)q^6 + 6(L-4)q^5r + 5(3L-14)q^4r^2 \right. \\
& + 20(L-5)q^3r^3 + 5(3L-14)q^2r^4 + 6(L-4)qr^5 + 2(L-4)r^6 \left. \right] \lambda'^3 \\
& - \frac{(q^2 + qr + r^2)}{16} \left[5(L-4)q^6 + 15(L-4)q^5r + (50L-247)q^4r^2 \right. \\
& + (75L-394)q^3r^3 + (50L-247)q^2r^4 \\
& \left. \left. + 15(L-4)qr^5 + 5(L-4)r^6 \right] \lambda'^4 + \dots \right\} \quad (4.2.14)
\end{aligned}$$

(for convenience in eventual comparison with the gauge theory, the eigenvalues have been expanded to $O(\lambda'^4)$). The corresponding result for the $SO(4)_{S^5}$ vector multiplet is

$$\begin{aligned}
\delta E_{S^5}(q, r, J, L) \approx & \frac{1}{J} \left\{ (L-8)(q^2 + qr + r^2)\lambda' - \frac{1}{2}(L-6)(q^2 + qr + r^2)^2\lambda'^2 \right. \\
& + \frac{3}{16} \left[2(L-6)q^6 + 6(L-6)q^5r + (15L-92)q^4r^2 \right. \\
& + 4(5L-31)q^3r^3 + (15L-92)q^2r^4 + 6(L-6)qr^5 + 2(L-6)r^6 \left. \right] \lambda'^3 \\
& - \frac{(q^2 + qr + r^2)}{16} \left[5(L-6)q^6 + 15(L-6)q^5r + (50L-309)q^4r^2 \right. \\
& + 3(25L-156)q^3r^3 + (50L-309)q^2r^4 \\
& \left. \left. + 15(L-6)qr^5 + 5(L-6)r^6 \right] \lambda'^4 + \dots \right\}. \quad (4.2.15)
\end{aligned}$$

Finally, the result for the spinor multiplet is

$$\begin{aligned}
\delta E_{\text{Fermi}}(q, r, J, L) \approx & \frac{1}{J} \left\{ (L-7)(q^2 + qr + r^2)\lambda' - \frac{1}{2}(L-5)(q^2 + qr + r^2)^2\lambda'^2 \right. \\
& + \frac{3}{16} \left[2(L-5)q^6 + 6(L-5)q^5r + 3(5L-27)q^4r^2 + 4(5L-28)q^3r^3 \right. \\
& \left. \left. + 3(5L-27)q^2r^4 + 6(L-5)qr^5 + 2(L-5)r^6 \right] \lambda'^3 \right.
\end{aligned}$$

$$\begin{aligned}
& -\frac{(q^2 + qr + r^2)}{16} \left[5(L-5)q^6 + 15(L-5)q^5r + 2(25L-139)q^4r^2 + (75L-431)q^3r^3 \right. \\
& \left. + 2(25L-139)q^2r^4 + 15(L-5)qr^5 + 5(L-5)r^6 \right] \lambda'^4 + \dots \Big\} . \quad (4.2.16)
\end{aligned}$$

It is important to remember that, to obtain the energies of the states as opposed to the energy shifts δE , we must add the BMN energy of the original degenerate multiplet to the above results:

$$\begin{aligned}
E_{\text{BMN}} &= \sqrt{1 + \lambda' q^2} + \sqrt{1 + \lambda' r^2} + \sqrt{1 + \lambda' (q+r)^2} \\
&= 3 + (q^2 + r^2 + qr)\lambda' - \frac{1}{4}(q^2 + r^2 + qr)^2 \lambda'^2 + \dots . \quad (4.2.17)
\end{aligned}$$

We can conclude from the above formulas that all three multiplets have a common internal level spacing given by the following function of λ' and mode indices:

$$\begin{aligned}
\frac{\delta E}{\delta L} &\approx \frac{1}{J} \Big\{ (q^2 + qr + r^2)\lambda' - \frac{1}{2}[(q^2 + qr + r^2)^2] \lambda'^2 \\
&+ \frac{3}{16} [2q^6 + 6q^5r + 15q^4r^2 + 20q^3r^3 + 15q^2r^4 + 6qr^5 + 2r^6] \lambda'^3 \\
&- \frac{(q^2 + qr + r^2)}{16} [5q^6 + 15q^5r + 50q^4r^2 + 75q^3r^3 \\
&+ 50q^2r^4 + 15qr^5 + 5(r^6)] \lambda'^4 + \dots \Big\} . \quad (4.2.18)
\end{aligned}$$

We have expanded in powers of λ' , but an all-orders formula can easily be constructed. The multiplets are displaced from one another by shifts that also depend on λ' and mode indices. We note that the one-loop degeneracy between different multiplets (see table 4.10) is preserved to second order in λ' , but is broken explicitly at three loops. At this order and beyond, each multiplet acquires a constant overall (L -independent) shift relative to the other two.

4.2.2 Two equal mode indices: ($q = r = n$, $s = -2n$)

An independent analysis is required when two mode indices are equal (specifically, we choose $q = r = n$, $s = -2n$). The all-loop matrix elements are complicated and we will

refrain from giving explicit expressions for them (though the complete formulas can be found at the following URL: <http://theory.caltech.edu/~swanson/MMA1/mma1.html>). As in the unequal mode index case, however, exact eigenvalues can easily be extracted by projection onto certain protected subsectors. In particular, the energy shift for states created by three bosonic mode creation operators with symmetric-traceless $SO(4)_{AdS}$ vector indices (the $\mathfrak{sl}(2)$ sector) turns out to be

$$\begin{aligned} \delta E_{AdS}(n, J) &= -\frac{n^2 \lambda'}{J(1 + n^2 \lambda') \sqrt{4n^2 + 1/\lambda'}} \left\{ \sqrt{4n^2 + \frac{1}{\lambda'}} (3 + 4n^2 \lambda') + \omega_n (4 + 8n^2 \lambda') \right\} \\ &\approx \frac{1}{J} \left\{ -7n^2 \lambda' + n^4 \lambda'^2 - \frac{17}{2} n^6 \lambda'^3 + \dots \right\}. \end{aligned} \quad (4.2.19)$$

The leading order term in the small- λ' expansion is the $-7/3$ [280_B] level $L = 4$ eigenvalue in the Λ_2 multiplet in table 4.13. The energy shift of the $SO(4)_{S^5}$ partners of these states (belonging to the $\mathfrak{su}(2)$ closed sector) is

$$\begin{aligned} \delta E_{S^5}(n, J) &= -\frac{n^2 \lambda'}{J(1 + n^2 \lambda') \sqrt{4n^2 + 1/\lambda'}} \left\{ \sqrt{4n^2 + \frac{1}{\lambda'}} (5 + 4n^2 \lambda') + \omega_n (6 + 8n^2 \lambda') \right\} \\ &\approx \frac{1}{J} \left\{ -11n^2 \lambda' + 8n^4 \lambda'^2 - \frac{101}{4} n^6 \lambda'^3 + \dots \right\}. \end{aligned} \quad (4.2.20)$$

The one-loop correction corresponds to the $-11/3$ [280_B] level in the Λ_1 submultiplet of table 4.13. As noted above, the protected symmetrized-fermion ($\mathfrak{su}(1|1)$) sector does not appear when two mode indices are equal. As in the previous section, we can do a numerical diagonalization of the full perturbation matrix to verify that the predicted eigenvalues are indeed exact and closed, but we will omit the details.

By invoking the angular momentum shift $J \rightarrow J + 2 - L/2$ in the BMN limit, we can use the energy shift of the $L = 4$ level to recover the exact energy shifts of all other levels in the superconformal multiplets of table 4.13. The energy shifts of the vector multiplet containing the protected $SO(4)_{AdS}$ bosonic irrep at level $L = 4$ are

given by the expression

$$\delta E_{AdS}(n, J, L) \approx \frac{1}{J} \left\{ \frac{1}{2}(3L - 19)n^2\lambda' - \frac{1}{2}(9L - 38)n^4\lambda'^2 + \frac{1}{8}(99L - 464)n^6\lambda'^3 - \frac{1}{16}(645L - 3160)n^8\lambda'^4 + \dots \right\}. \quad (4.2.21)$$

The shifts of the multiplet containing the protected $SO(4)_{S^5}$ bosonic irrep are given by

$$\delta E_{S^5}(n, J, L) \approx \frac{1}{J} \left\{ \frac{1}{2}(3L - 23)n^2\lambda' - \frac{1}{2}(9L - 52)n^4\lambda'^2 + \frac{1}{8}(99L - 598)n^6\lambda'^3 - \frac{1}{16}(645L - 3962)n^8\lambda'^4 + \dots \right\}. \quad (4.2.22)$$

Once again, we note that in order to get energies, rather than energy shifts, one must append the BMN energy of the original degenerate multiplet to these results. Unlike the unequal mode index case, there is no accidental degeneracy between superconformal multiplets spanning the three-impurity space, even at one loop in λ' . The level spacings within the two superconformal multiplets are the same, but the multiplets are offset from each other by an L -independent shift (but one that depends on λ' and mode indices).

4.3 Gauge theory anomalous dimensions

In the previous sections, we have given a complete analysis of the perturbed energy spectrum of three-impurity string states. The “data” are internally consistent in the sense that the perturbed energy levels organize themselves into proper superconformal multiplets of the classical nonlinear sigma model governing the string worldsheet dynamics. Since the quantization procedure leaves only a subgroup of the full symmetry group as a manifest, linearly realized symmetry, this is by itself a nontrivial check on the consistency of the action and quantization procedure. To address the issue of AdS/CFT duality, we must go further and compare the string energy spectrum

with the anomalous dimensions of gauge theory operators dual to the three-impurity string states.

As discussed in previous chapters, the task of finding the anomalous dimensions of BMN operators in the limit of large R -charge and dimension D , but finite $\Delta = D - R$, is greatly simplified by the existence of an equivalence between the dilatation operator of $\mathcal{N} = 4$ SYM and the Hamiltonian of a one-dimensional spin chain. The one-loop spin chain Hamiltonian has only nearest-neighbor interactions (in the planar large- N_c limit) and is of limited complexity. This is tempered by the fact that the higher-loop gauge theory physics is encoded in increasingly long-range spin chain interactions that generate a rapidly growing number of possible terms in the Hamiltonian [31]. Fixing the coefficients of all these terms by comparison with diagrammatic computations would be a very impractical approach. Fortunately, Beisert was able to show that, at least for BMN operators in the $\mathfrak{su}(2)$ closed subsector, general requirements (such as the existence of a well-defined BMN scaling limit) suffice to fix the form of the spin chain Hamiltonian out to three-loop order [30, 33]. In Chapter 2 we discussed the use of these higher-loop spin chains to generate the information we need on the anomalous dimensions of three-impurity operators: we will rely on these results for our comparison with the three-impurity string theory predictions computed above.

We have already noted that there are three closed subsectors of BMN operators in which impurities taken from a subalgebra of the full superconformal algebra mix only with themselves: we have referred to them as the $\mathfrak{sl}(2)$, $\mathfrak{su}(2)$ (both bosonic) and $\mathfrak{su}(1|1)$ (fermionic) sectors. We will focus our attention on these sectors because their spin chain description is simple and their anomalous dimensions fix the dimensions of the remaining three-impurity operators in the theory. Spin chain Hamiltonians incorporating higher-loop-order gauge theory physics have been constructed for the $\mathfrak{su}(2)$ and $\mathfrak{su}(1|1)$ sectors but, as far as we know, the $\mathfrak{sl}(2)$ spin chain is known only to one-loop order.⁴

Although these spin chains are integrable, methods such as the Bethe ansatz tech-

⁴See ref. [52] for important progress on this problem.

nique do not immediately yield the desired results for all multiple-impurity anomalous dimensions of interest. Minahan and Zarembo did use the Bethe ansatz for the one-loop $\mathfrak{so}(6)$ spin chain (of which the exactly closed $\mathfrak{su}(2)$ system is a subsector) to obtain approximate multi-impurity anomalous dimensions [39], but we need results for all sectors and for higher-loop spin chains. As mentioned above, the $\mathfrak{sl}(2)$ spin chain has phenomenological applications and has been extensively developed in that context. It is therefore possible that some of the results we need can be extracted from the relevant literature.⁵ In the end, since we are looking for a unified approach that can handle all sectors and any number of loops, we decided that numerical methods are, for the present purposes, an effective way to extract the information we need about gauge theory anomalous dimensions. Since Bethe ansatz equations exist for most of the results that are of interest to us, the numerical results obtained here can be checked against the Bethe-ansatz methodology: these exercises were performed in Chapter 2 above (see ref. [95]).

We begin with a discussion of the bosonic $\mathfrak{sl}(2)$ sector. Recall from Chapter 2 that for total R -charge K (the R -charge is equal to the number of lattice sites K in this sector), the basis for this system consists of single-trace operators of the form

$$\mathrm{Tr}(\nabla^I Z Z^{K-1}), \mathrm{Tr}(\nabla^{I-1} Z \nabla Z Z^{K-2}), \mathrm{Tr}(\nabla^{I-1} Z Z \nabla Z Z^{K-3}), \dots, \quad (4.3.1)$$

where Z is the $SO(6)$ Yang-Mills boson carrying one unit of R -charge, ∇ is a space-time covariant derivative operator that scales under the chosen $\mathfrak{sl}(2)$ subgroup of the Lorentz group ($\nabla \equiv \nabla_1 + i\nabla_2$), I is the total impurity number and the full basis contains all possible distributions of ∇ operators among the Z fields. Conservation of various $U(1)$ subgroups of the R -symmetry group ensures that operators of this type mix only among themselves to all orders in the gauge theory (as long as we work in the 't Hooft large- N_c limit). This gauge theory closed subsector corresponds to the symmetric traceless irrep of $SO(4)_{AdS}$ bosons in the string theory (states whose

⁵We thank A. Belitsky for making us aware of this literature and for helpful discussions on this point.

energy shifts are given in eqns. (4.2.9) and (4.2.19)).

To compare the results of Chapter 2 (see eqn. (2.3.7)) with the string theory predictions of eqns. (4.2.9) and (4.2.19), we reorganize those results as follows: we reinstate the BMN energy of the degenerate multiplet (4.2.17) (expanded to first order in λ'); we replace λ' with λ/J^2 and replace J by K . This gives specific string theory predictions for the large- K scaling of one-loop anomalous dimensions of the AdS closed sector. As usual, there are two distinct cases: for unequal mode indices ($q \neq r \neq s = -q - r$), we have

$$E_{AdS}(q, r, K) = 3 + (K - 2)(q^2 + r^2 + qr) \frac{\lambda}{K^3} + O(K^{-4}) . \quad (4.3.2)$$

(Note that here we label mode indices with q, r, s instead of the k_1, k_2, k_3 triplet used in Chapter 2.) For pairwise equal mode indices ($n, n, -2n$) we have

$$E_{AdS}(n, K) = 3 + (3K - 7)n^2 \frac{\lambda}{K^3} + O(K^{-4}) . \quad (4.3.3)$$

For convenience, we redisplay the numerical gauge theory predictions from the $\mathfrak{sl}(2)$ spin chain from eqn. (2.3.7):

$$\begin{aligned} E_{\mathfrak{sl}(2)}^{(1,2)} &= (k_1^2 + k_1 k_2 + k_2^2) , & E_{\mathfrak{sl}(2)}^{(1,3)} / E_{\mathfrak{sl}(2)}^{(1,2)} &= -2 , & k_1 &\neq k_2 \neq k_3 , \\ E_{\mathfrak{sl}(2)}^{(1,2)} &= 3n^2 , & E_{\mathfrak{sl}(2)}^{(1,3)} / E_{\mathfrak{sl}(2)}^{(1,2)} &= -7/3 , & k_1 &= k_2 = n, \, k_3 = -2n . \end{aligned} \quad (4.3.4)$$

The string predictions in eqns. (4.3.2) and (4.3.3) match the expected virial scaling of the spin chain eigenvalues displayed in eqn. (2.3.7), with the specific identifications

$$E_{AdS}^{(1,2)} = (q^2 + r^2 + qr) , \quad E_{AdS}^{(1,3)} = -2(q^2 + r^2 + qr) , \quad E_{AdS}^{(1,3)} / E_{AdS}^{(1,2)} = -2 \quad (4.3.5)$$

for $q \neq r \neq s = -q - r$, or

$$E_{AdS}^{(1,2)} = 3n^2, \quad E_{AdS}^{(1,3)} = -7n^2, \quad E_{AdS}^{(1,3)}/E_{AdS}^{(1,2)} = -7/3 \quad (4.3.6)$$

for $q = r = n$ and $s = -2n$.

At this point it is appropriate to say a few words about the role of integrability in this problem. It was first argued in [59] that the complete GS action of IIB superstring theory on $AdS_5 \times S^5$ is integrable. Integrability has since taken a central role in studies of the AdS/CFT correspondence, as any precise non-perturbative understanding of integrability on both sides of the duality would be extremely powerful. Integrability on either side of the duality gives rise to an infinite tower of hidden charges that can be loosely classified as either local (Abelian) or non-local (non-Abelian). In the Abelian sector, contact between the integrable structures of gauge theory and semiclassical string theory (a subject that was first investigated in [41]) has been made to two loops in λ (see, e.g., [29, 45, 50, 96]). (The corresponding problem in the non-local sector was addressed to one-loop order in [47, 48].) One of the local gauge theory charges can be shown to anticommute in the $\mathfrak{su}(2)$ sector with a parity operator P (to three loops in λ), whose action on a single-trace state in the gauge theory is to invert the order of all fields within the trace [31, 33]. Furthermore, this operator can be shown to connect states of opposite parity. These facts imply that all eigenstates in the spectrum connected by P must be degenerate. These degenerate states are known as parity pairs and their existence can be interpreted as a necessary (but not sufficient) condition for integrability. The spectrum in table 2.8 from Chapter 2 exhibits such a degeneracy and makes it clear that parity pairs are simply distinct states whose lattice momenta (or worldsheet mode indices) are related by an overall sign flip. Since the net momentum of allowed states is zero, parity pair states can in principle scatter into each other, and their degeneracy is a nontrivial constraint on the interactions. As a small caveat, we note that lattice momentum conservation implies that mixing of parity-pair states can only occur via connected three-body (or higher)

interactions. As the virial analysis shows, at the order to which we are working, only two-body interactions are present and the parity pair degeneracy is automatic. The same remark applies to the string theory analysis to $O(J^{-1})$ in the curvature expansion. A calculation of the string theory spectrum carried out to $O(J^{-2})$ is needed to see whether parity pair degeneracy survives string worldsheet interactions; further discussion of this point will be given in Chapter 6 (see also ref. [71]).

We now turn to the closed $\mathfrak{su}(2)$ sector of gauge theory operators, corresponding to the symmetric-traceless bosonic $SO(4)_{S^5}$ sector of the string theory. The operator basis for this sector consists of single-trace monomials built out of two complex scalar fields Z and ϕ , where Z is the complex scalar carrying one unit of charge under the $U(1)$ R -charge subgroup and ϕ is one of the two scalars with zero R -charge, transforming as an $SO(4)$ vector in the $SO(6) \simeq U(1)_R \times SO(4)$ decomposition of the full R -symmetry group of the gauge theory. The collection of operators

$$\text{tr}(\phi^I Z^{K-I}), \text{tr}(\phi^{I-1} Z \phi Z^{K-I-1}), \text{tr}(\phi^{I-2} Z \phi^2 Z^{K-I-1}), \dots \quad (4.3.7)$$

(and all possible permutations, modulo cyclic equivalence, of the K factors) forms a basis with I impurities and R -charge equal to $K - I$. The anomalous dimension operator simply permutes these monomials among themselves in ways that get more elaborate as we go to higher loop orders in the gauge theory. The relevant gauge theory predictions from Chapter 2 are given at one-loop order in eqn. (2.1.28) and at two-loop order in eqn. (2.1.42). For convenience, we reproduce those results here:

$$\begin{aligned} E_{\mathfrak{su}(2)}^{(1,2)} &= (k_1^2 + k_2^2 + k_3^2)/2, & k_1 + k_2 + k_3 &= 0, \\ E_{\mathfrak{su}(2)}^{(1,3)}/E_{\mathfrak{su}(2)}^{(1,2)} &= 2, & (k_1 \neq k_2 \neq k_3), \\ E_{\mathfrak{su}(2)}^{(1,3)}/E_{\mathfrak{su}(2)}^{(1,2)} &= \frac{7}{3}, & (k_1 = k_2, k_3 = -2k_1), \\ E_{\mathfrak{su}(2)}^{(2,4)} &= -(k_1^2 + k_2^2 + k_3^2)^2/16, & k_1 + k_2 + k_3 &= 0, \\ E_{\mathfrak{su}(2)}^{(2,5)}/E_{\mathfrak{su}(2)}^{(2,3)} &= 8, & (k_1 \neq k_2 \neq k_3), \\ E_{\mathfrak{su}(2)}^{(2,5)}/E_{\mathfrak{su}(2)}^{(2,3)} &= \frac{76}{9}, & (k_1 = k_2, k_3 = -2k_1). \end{aligned} \quad (4.3.8)$$

To compare with string theory results for the bosonic symmetric-traceless $SO(4)_{S^5}$ sector eigenvalues, we need to recast eqns. (4.2.10) and (4.2.20) as expansions in powers of λ and K^{-1} . We denote by $E_{S^5}^{(n,m)}$ the coefficient of $\lambda^n K^{-m}$ in the large- K expansion of the string theory energies: they can be directly compared with the corresponding quantities extracted from the numerical spin chain analysis. The string theory predictions for scaling coefficients, up to second order in λ , are given in table 4.16. As usual, the predictions for three-impurity states with unequal mode indices

$E_{S^5}^{(n,m)}$	$(q \neq r \neq s)$	$(q = r = n)$
$E_{S^5}^{(1,2)}$	$(q^2 + qr + r^2)$	$3n^2$
$E_{S^5}^{(1,3)}$	$2(q^2 + qr + r^2)$	$7n^2$
$E_{S^5}^{(2,4)}$	$-\frac{1}{4}(q^2 + qr + r^2)^2$	$-\frac{9}{4}n^4$
$E_{S^5}^{(2,5)}$	$-2(q^2 + qr + r^2)^2$	$-19n^4$

Table 4.16: String predictions for $\mathfrak{su}(2)$ scaling coefficients, to two loops

have to be stated separately from those for states with two equal mode indices. We take these results as strong evidence that the string theory analysis agrees with the gauge theory up to $O(\lambda^2)$ in this sector.

We now turn to a discussion of gauge theory physics beyond two loops. As it happens, the three-loop Hamiltonian can be fixed up to two unknown coefficients (α_1 and α_2) by basic field theory considerations [30]:

$$\begin{aligned}
H_6^{\mathfrak{so}(6)} = & (60 + 6\alpha_1 - 56\alpha_2) \{ \} + (-104 + 14\alpha_1 + 96\alpha_2) \{ 0 \} \\
& + (24 + 2\alpha_1 - 24\alpha_2) (\{ 0, 1 \} + \{ 1, 0 \}) + (4 + 6\alpha_1) \{ 0, 2 \} \\
& (-4 + 4\alpha_2) (\{ 0, 1, 2 \} + \{ 2, 1, 0 \}) - \alpha_1 (\{ 0, 2, 1 \} + \{ 1, 0, 2 \}) . \quad (4.3.9)
\end{aligned}$$

Originally, these coefficients were determined by demanding proper BMN scaling in the theory and that the dynamics be integrable at three loops; these assumptions set $\alpha_{1,2} = 0$. By studying an $\mathfrak{su}(2|3)$ spin chain model, Beisert [33] was subsequently able to show that independent symmetry arguments, along with BMN scaling, uniquely

set $\alpha_1 = \alpha_2 = 0$ (thus proving integrability at three loops).

As described in Chapter 2, the three-loop Hamiltonian $H_6^{\mathfrak{su}(2)}$ can be treated as a second-order correction to $H_2^{\mathfrak{su}(2)}$. This allows us to numerically evaluate the $O(\lambda^3)$ contribution to the spectrum by using second-order Rayleigh-Schrödinger perturbation theory (there is an intermediate state sum involved, but since we are doing the calculation numerically, this is not a serious problem). There is also the issue of degeneracy but the existence of a higher conserved charge once again renders the problem effectively non-degenerate. The resulting three-loop data for large- K was fit in Chapter 2 to a power series in K^{-1} to read off the expansion coefficients $E_{\mathfrak{su}(2)}^{3,n}$. It turns out that, to numerical precision, the coefficients are non-vanishing only for $n > 5$ (as required by BMN scaling). The results of this program are reproduced for convenience from Chapter 2 in table 4.17, where they are compared with string theory predictions derived (in the manner described in previous paragraphs) from eqn. (4.2.10). (The accuracy of the match is displayed in the last column of table 4.17.) The important point is that there is substantial disagreement with string results at $O(\lambda^3)$ for all energy levels: the low-lying states exhibit a mismatch ranging from roughly 19% to 34%, and there is no evidence that this can be repaired by taking data on a larger range of lattice sizes. There is apparently a general breakdown of the correspondence between string theory and gauge theory anomalous dimensions at three loops, despite the precise and impressive agreement at first and second order. This disagreement was first demonstrated in the two-impurity regime [26]. It is perhaps not surprising that the three-loop disagreement is reproduced in the three-impurity regime, but it provides us with more information that may help to clarify this puzzling phenomenon.

The same exercise can be repeated for the closed $\mathfrak{su}(1|1)$ fermionic sector, whose string theory dual is comprised of pure fermionic states symmetrized in $SO(4) \times SO(4)$ indices in either the $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ or $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ irreps (projected onto Π_{\pm} subspaces). The spin chain system is embedded in Beisert's $\mathfrak{su}(2|3)$ model, where the fermionic sector of the Hamiltonian has been recorded up to two-loop order [33]. Since the

$E_{\mathfrak{su}(2)}^{(3,7)}/E_{\mathfrak{su}(2)}^{(3,6)}$	String Modes (q, r, s)	%Error
16.003	$(1, 0, -1)$	33%
14.07	$(1, 1, -2)$	19%
14.07	$(-1, -1, 2)$	19%
16.03	$(2, 0, -2)$	34%
14.37	$(1, 2, -3)$	22%
14.37	$(-1, -2, 3)$	22%
15.96	$(3, 0, -3)$	30%

Table 4.17: Three-loop numerical spectrum of three-impurity $\mathfrak{su}(2)$ subsector and mismatch with string predictions

relevant points of the numerical gauge/string comparison have already been made, we will simply state the one- and two-loop results. (The large- K spectrum of the three-loop contribution is scrutinized in Chapter 2; one again finds disagreement with string theory.)

In this sector, the R -charge and the lattice length are related by $J = K - I/2$. The fermionic one- and two-loop string predictions are therefore found from eqn. (4.2.11) to be

$$\begin{aligned}
E_{\text{Fermi}}^{(1,2)} &= (q^2 + qr + r^2) , & E_{\text{Fermi}}^{(1,3)} &= 0 \\
E_{\text{Fermi}}^{(2,4)} &= -\frac{1}{4}(q^2 + qr + r^2)^2 , & E_{\text{Fermi}}^{(2,5)} &= -(q^2 + qr + r^2)^2 .
\end{aligned} \quad (4.3.10)$$

As noted above, this sector does not admit states with equivalent mode indices.

Reproducing the results from the gauge theory analysis in Chapter 2, we find precise agreement:

$$\begin{aligned}
E_{\mathfrak{su}(1|1)}^{(1,2)} &= (k_1^2 + k_1 k_2 + k_2^2) , & E_{\mathfrak{su}(1|1)}^{(1,3)} &= 0 , \\
E_{\mathfrak{su}(1|1)}^{(2,4)} &= -\frac{1}{4}(k_1^2 + k_1 k_2 + k_2^2)^2 , & E_{\mathfrak{su}(1|1)}^{(2,5)} &= -(k_1^2 + k_1 k_2 + k_2^2)^2 .
\end{aligned} \quad (4.3.11)$$

The two-loop data are obtained using the same first-order perturbation theory treatment described above in the $\mathfrak{su}(2)$ sector (the results are recorded in table 2.6 above).

The two-loop spectrum is subject to stronger K^{-1} corrections, but the data are still convincing and could be improved by running the extrapolation out to larger lattice sizes. Nonetheless, the close agreement for the low-lying levels corroborates the match between gauge and string theory up to two-loop order.

4.4 Discussion

The BMN/pp-wave mechanism has emerged as a useful proving ground for the postulates of the AdS/CFT correspondence. When the full Penrose limit is lifted, a rich landscape emerges, even in the two-impurity regime, upon which the string and gauge theory sides of the duality have exhibited an intricate and impressive match to two loops in the gauge coupling and first nontrivial order in the curvature expansion. While the conditions under which agreement is obtained are substantially more demanding in the higher-impurity problem, we have shown that this agreement is maintained for three-impurity string states and SYM operators. Although the two-loop agreement survives at the three-impurity level, we have also confirmed the previously observed mismatch at three loops in the gauge theory coupling. In the end, the analyses carried out here will provide an extremely stringent test of any proposed solution to this vexing problem.

Chapter 5

N impurities

In Chapters 3 and 4 we analyzed the first curvature correction to the spectrum of string states in the pp-wave limit of $AdS_5 \times S^5$. The string energies in this setting correspond in the gauge theory to the difference between operator scaling dimensions and R -charge ($\Delta \equiv D - R$), and states are arranged into superconformal multiplets according to the $\mathfrak{psu}(2, 2|4)$ symmetry of the theory. The fully supersymmetric two-excitation (or two-impurity) system, for example, is characterized by a 256-dimensional supermultiplet of states built on a scalar primary. The complete spectrum of this system was successfully matched to corresponding SYM operator dimensions in Chapter 3 to two loops in the modified 't Hooft coupling $\lambda' = \lambda/J^2$ (see also [26, 27]). A three-loop mismatch between the gauge and string theory results discovered therein comprises a long-standing and open problem in these studies, one which has appeared in several different contexts (see, e.g., [29, 34, 37]). This was extended to the three-impurity, 4,096-dimensional supermultiplet of string states in Chapter 4 (see also [38]), where precise agreement with the corresponding gauge theory was again found to two-loop order, and a general disagreement reappeared at three loops. In the latter study, three-impurity string predictions were compared with corresponding gauge theory results derived both from the virial technique described in Chapter 2 and the long-range Bethe ansatz of [32] (which overlaps at one loop with the original $\mathfrak{so}(6)$ system studied in [39]).

In the present chapter we generalize the string side of these investigations by

computing, directly from the Hamiltonian, various N -impurity spectra of IIB superstring theory at $O(J^{-1})$ in the large- J curvature expansion near the pp-wave limit of $AdS_5 \times S^5$. We focus on the bosonic $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ sectors, which are characterized by N symmetric-traceless bosonic string excitations in the S^5 and AdS_5 subspaces, respectively. Based on calculations in these sectors, we also formulate a conjecture for the N -impurity spectrum of states in a protected $\mathfrak{su}(1|1)$ sector composed of N fermionic excitations symmetrized in their $SO(4) \times SO(4)$ spinor indices. We then describe the complete supermultiplet decomposition of the N -impurity spectrum to two loops in λ' using a simple generalization of the two- and three-impurity cases.

We note here that a new Bethe ansatz for the string theory has been proposed by Arutyunov, Frolov and Staudacher [44] that is meant to diagonalize the fully quantized string sigma model in the $\mathfrak{su}(2)$ sector to all orders in $1/J$ and λ' (see the discussion in Chapter 1). This ansatz was shown in [44] to reproduce the two- and three-impurity spectra of quantized string states near the pp-wave limit detailed in [27, 38]. The methods developed here allow us to check their formulas directly against the string theory for any impurity number at $O(J^{-1})$, and we find that our general $\mathfrak{su}(2)$ string eigenvalues agree to all orders in λ' with their $\mathfrak{su}(2)$ string Bethe ansatz! We compute the N -impurity energy spectra of the $\mathfrak{su}(2)$, $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ closed sectors of this system in Section 5.1, and generalize the complete N -impurity supermultiplet structure of the theory to two-loop order in λ' in Section 5.2.

5.1 N -impurity string energy spectra

As described above, our string vacuum state carries the S^5 string angular momentum J and is labeled by $|J\rangle$; the complete Fock space of string states is generated by acting on $|J\rangle$ with any number of the creation operators $a_n^{A\dagger}$ (bosonic) and $b_n^{\alpha\dagger}$ (fermionic), where the lower indices n, m, l, \dots denote mode numbers. The excitation number of string states (defined by the number of creation oscillators acting on the ground state) will also be referred to as the impurity number, and string states with a total

of $N_B + N_F = N$ impurities will contain N_B bosonic and N_F fermionic impurities:

$$|N_B, N_F; J\rangle \equiv \underbrace{a_{n_1}^{A_1\dagger} a_{n_2}^{A_2\dagger} \dots a_{n_{N_B}}^{A_{N_B}\dagger}}_{N_B} \underbrace{b_{n_1}^{\alpha_1\dagger} b_{n_2}^{\alpha_2\dagger} \dots b_{n_{N_F}}^{\alpha_{N_F}\dagger}}_{N_F} |J\rangle . \quad (5.1.1)$$

States constructed in this manner fall into two disjoint subsectors populated by space-time bosons (N_F even) and spacetime fermions (N_F odd). In this notation the pure-boson states $|N_B, 0; J\rangle$ are mixed only by H_{BB} and the pure-fermion states $|0, N_F; J\rangle$ are acted on by H_{FF} . The more general spacetime-boson states $|N_B, \text{even}; J\rangle$ are acted on by the complete interaction Hamiltonian H_{int} , as are the spacetime-fermion states $|N_B, \text{odd}; J\rangle$. There is of course no mixing between spacetime bosons and fermions; this block-diagonalization is given schematically in table 5.1.

H_{int}	$ N_B, 0; J\rangle$	$ N_B, \text{even}; J\rangle$	$ N_B, \text{odd}; J\rangle$	$ 0, \text{odd}; J\rangle$
$\langle N_B, 0; J $	H_{BB}	H_{BF}		
$\langle N_B, \text{even}; J $	H_{BF}	$H_{BB} + H_{BF} + H_{FF}$		
$\langle N_B, \text{odd}; J $			$H_{BB} + H_{BF} + H_{FF}$	H_{BF}
$\langle 0, \text{odd}; J $			H_{BF}	H_{FF}

Table 5.1: Interaction Hamiltonian on N -impurity string states ($N_B + N_F = N$)

The full interaction Hamiltonian can be further block-diagonalized by projecting onto certain protected sectors of string states, and we will focus in this study on three such sectors. Two of these sectors are spanned by purely bosonic states $|N_B, 0; J\rangle$ projected onto symmetric-traceless irreps in either the $SO(4)_{AdS}$ or $SO(4)_{S^5}$ subspaces. Another sector that is known to decouple at all orders in λ' is comprised of purely fermionic states $|0, N_F; J\rangle$ projected onto either of two subspaces of $SO(4) \times SO(4)$ labeled, in an $SU(2)^2 \times SU(2)^2$ notation, by $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ and $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$, and symmetrized in spinor indices. Each of these sectors can also be labeled by the subalgebra of the full superconformal algebra that corresponds to the symmetry under which they are invariant. The bosonic $SO(4)_{AdS}$ and $SO(4)_{S^5}$ sectors are labeled by $\mathfrak{sl}(2)$ and $\mathfrak{su}(2)$ subalgebras, respectively, while the two fermionic sectors fall into $\mathfrak{su}(1|1)$ subsectors

of the closed $\mathfrak{su}(2|3)$ system studied in [32, 33, 69].

In the large- J expansion about the free pp-wave theory, we will isolate $O(J^{-1})$ corrections to the energy eigenvalues of N -impurity string states according to

$$E(\{q_j\}, N, J) = \sum_{j=1}^N \sqrt{1 + q_j^2 \lambda'} + \delta E(\{q_j\}, N, J) + O(J^{-2}) . \quad (5.1.2)$$

The spectrum is generically dependent upon λ' , J and the mode numbers $\{n_j\}, \{q_j\}, \dots$, where j is understood to label either the complete set of impurities ($j = 1, \dots, N$) or some subset thereof (e.g., $j = 1, \dots, N_F$). The leading order term in this expansion is the N -impurity free energy of states on the pp-wave geometry, and $\delta E(\{q_j\}, N, J)$ always enters at $O(J^{-1})$. When it becomes necessary, we will also expand the $O(1/J)$ energy shift in the small- λ' loop expansion:

$$\delta E(\{q_j\}, N, J) = \sum_{i=1}^{\infty} \delta E^{(i)}(\{q_j\}, N, J) (\lambda')^i . \quad (5.1.3)$$

Finding the explicit form of $\delta E(\{q_j\}, N, J)$ for N -impurity string states in certain interesting sectors of the theory will be our primary goal. As a side result, however, we will see that the spectrum of *all* states in the theory will be determined to two-loop order in λ' by the specific eigenvalues we intend to compute.

We begin by noting that the canonical commutation relations of the bosonic fields x^A and p_A allow us to expand H_{BB} in bosonic creation and annihilation operators using

$$\begin{aligned} x^A(\sigma, \tau) &= \sum_{n=-\infty}^{\infty} x_n^A(\tau) e^{-ik_n \sigma} , \\ x_n^A(\tau) &= \frac{i}{\sqrt{2\omega_n}} \left(a_n^A e^{-i\omega_n \tau} - a_{-n}^{A\dagger} e^{i\omega_n \tau} \right) , \end{aligned} \quad (5.1.4)$$

where $k_n = n$ are integer-valued, $\omega_n = \sqrt{p_-^2 + n^2}$ and the operators a_n^A and $a_n^{A\dagger}$ obey the usual relation $[a_m^A, a_n^{B\dagger}] = \delta_{mn} \delta^{AB}$. Since we are only interested in computing diagonal matrix elements of H_{BB} between physical string states with equal numbers

of excitations, we can restrict the oscillator expansion to contain only equal numbers of creation and annihilation operators (all other combinations automatically annihilate between equal-impurity string states). Explicitly, we obtain the following expansion from the results in Chapter 3:

$$\begin{aligned}
H_{\text{BB}} = & -\frac{1}{32p_-R^2} \sum \frac{\delta(n+m+l+p)}{\xi} \\
& \times \left\{ 2 \left[\xi^2 - (p_-^4 - k_l k_p k_n k_m) + \omega_n \omega_m k_l k_p + \omega_l \omega_p k_n k_m + 2\omega_n \omega_l k_m k_p \right. \right. \\
& + 2\omega_m \omega_p k_n k_l \left. \right] a_{-n}^{\dagger A} a_{-m}^{\dagger A} a_l^B a_p^B + 4 \left[\xi^2 - (p_-^4 - k_l k_p k_n k_m) - 2\omega_n \omega_m k_l k_p + \omega_l \omega_m k_n k_p \right. \\
& - \omega_n \omega_l k_m k_p - \omega_m \omega_p k_n k_l + \omega_n \omega_p k_m k_l \left. \right] a_{-n}^{\dagger A} a_{-l}^{\dagger B} a_m^A a_p^B + 2 \left[8k_l k_p a_{-n}^{\dagger i} a_{-l}^{\dagger j} a_m^i a_p^j \right. \\
& + 2(k_l k_p + k_n k_m) a_{-n}^{\dagger i} a_{-m}^{\dagger i} a_l^j a_p^j + (\omega_l \omega_p + k_l k_p - \omega_n \omega_m - k_n k_m) a_{-n}^{\dagger i} a_{-m}^{\dagger i} a_l^{j'} a_p^{j'} \\
& \left. \left. - 4(\omega_l \omega_p - k_l k_p) a_{-n}^{\dagger i} a_{-l}^{\dagger j'} a_m^i a_p^{j'} - (i, j \rightleftharpoons i', j') \right] \right\} , \tag{5.1.5}
\end{aligned}$$

where $\xi \equiv \sqrt{\omega_n \omega_m \omega_l \omega_p}$.

5.1.1 The $SO(4)_{S^5}$ ($\mathfrak{su}(2)$) sector

We begin in the $\mathfrak{su}(2)$ sector spanned by symmetric-traceless pure-boson states excited in the S^5 subspace. Because we are restricting our attention to $SO(4)_{S^5}$ states symmetric in their vector indices, we form the following oscillators:

$$a_n = \frac{1}{\sqrt{2}} (a_n^5 + i a_n^6) , \quad \bar{a}_n = \frac{1}{\sqrt{2}} (a_n^5 - i a_n^6) . \tag{5.1.6}$$

By taking matrix elements of the form

$$\langle J | a_{n_1} a_{n_2} \dots a_{n_{N_B}} (H_{\text{BB}}) a_{n_1}^{\dagger} a_{n_2}^{\dagger} \dots a_{n_{N_B}}^{\dagger} | J \rangle , \tag{5.1.7}$$

we can therefore select out excitations in the $(5, 6)$ -plane of the S^5 subspace and make the symmetric-traceless projection manifest. (More generally we can project onto any (n, m) -plane, as long as $n \neq m$ and both are chosen to lie in the S^5 subspace.)

There are two basic oscillator structures of H_{BB} in eqn. (5.1.5): one in which the creation (annihilation) operators are contracted in their $SO(4) \times SO(4)$ indices

$$a_{-n}^{\dagger A} a_{-m}^{\dagger A} a_l^B a_p^B ,$$

and one where pairs of creation and annihilation operators are contracted

$$a_{-n}^{\dagger A} a_{-l}^{\dagger B} a_m^A a_p^B .$$

In terms of the a_n and \bar{a}_n fields of eqn. (5.1.6), the former structure contains

$$a_{-n}^{\dagger A} a_{-m}^{\dagger A} a_l^B a_p^B \Big|_{(5,6)} = (a_{-n}^{\dagger} \bar{a}_{-m}^{\dagger} + \bar{a}_{-n}^{\dagger} a_{-m}^{\dagger}) (a_l \bar{a}_p + \bar{a}_l a_p) , \quad (5.1.8)$$

which cannot contribute to $\mathfrak{su}(2)$ matrix elements of the form appearing in (5.1.7).

The latter structure, however, contains

$$a_{-n}^{\dagger A} a_{-l}^{\dagger B} a_m^A a_p^B \Big|_{(5,6)} = \bar{a}_{-n}^{\dagger} \bar{a}_{-l}^{\dagger} \bar{a}_m \bar{a}_p + a_{-n}^{\dagger} a_{-l}^{\dagger} a_m a_p , \quad (5.1.9)$$

which will contribute to the $\mathfrak{su}(2)$ energy spectrum.

The string states appearing in the matrix element of eqn. (5.1.7) have been written in the generic form

$$a_{n_1}^{\dagger} a_{n_2}^{\dagger} \dots a_{n_{N_B}}^{\dagger} |J\rangle ,$$

and, as usual, they are subject to the level-matching condition

$$\sum_{j=1}^{N_B} n_j = 0 . \quad (5.1.10)$$

The complete set of mode indices $\{n_1, n_2, \dots, n_{N_B}\}$ can contain one or more subsets of indices that are equal, while still satisfying eqn. (5.1.10); this scenario complicates the calculation of energy eigenvalues to some extent. We will eventually compute

the eigenvalues of interest for completely general string states, but for purposes of illustration and to introduce our strategy we will start with the simplest case in which no two mode numbers are equal ($n_1 \neq n_2 \neq \dots \neq n_{N_B}$). To organize the presentation of this chapter, we will use mode numbers labeled by $\{n_j\}$ to denote those that are inequivalent from each other, while $\{q_j\}$ will be allowed to overlap. Between states with completely distinct mode indices, the oscillator structure in eqn. (5.1.9) exhibits the following matrix element:

$$\begin{aligned} \langle J | a_{n_1} a_{n_2} \dots a_{N_B} (a_{-n}^\dagger a_{-l}^\dagger a_m a_p) a_{n_1}^\dagger a_{n_2}^\dagger \dots a_{N_B}^\dagger | J \rangle \\ = \frac{1}{2} \sum_{\substack{j,k=1 \\ j \neq k}}^{N_B} \left(\delta_{n_j+n} \delta_{n_k+l} \delta_{n_j-m} \delta_{n_k-p} + \delta_{n_j+n} \delta_{n_k+l} \delta_{n_k-m} \delta_{n_j-p} \right. \\ \left. + \delta_{n_j+l} \delta_{n_k+n} \delta_{n_j-m} \delta_{n_k-p} + \delta_{n_j+l} \delta_{n_k+n} \delta_{n_k-m} \delta_{n_j-p} \right) . \end{aligned} \quad (5.1.11)$$

With this in hand, it is a straightforward exercise to compute the energy eigenvalue of the $SO(4)_{S^5}$ bosonic interaction Hamiltonian in the N_B -impurity symmetric-traceless irrep (with unequal mode indices): we simply attach the H_{BB} coefficient of the oscillator structure $a_{-n}^\dagger a_{-l}^\dagger a_m a_p$ to the right-hand side of eqn. (5.1.11) and carry out the summation over mode numbers. The result is remarkably compact:

$$\delta E_{S^5}(\{n_i\}, N_B, J) = -\frac{1}{J} \sum_{\substack{j,k=1 \\ j \neq k}}^{N_B} \frac{1}{2\omega_{n_j}\omega_{n_k}} \left[n_k^2 + n_j^2 (1 + n_k^2 \lambda') + n_j n_k (1 - \omega_{n_j} \omega_{n_k} \lambda') \right] . \quad (5.1.12)$$

This $\mathfrak{su}(2)$ formula can be checked against previously obtained string theory results in the two- and three-impurity regimes. Namely, the two-impurity eigenvalue computed above (and in [26, 27]) takes the form (which is exact in λ')

$$\delta E_{S^5}(n_1, n_2, J) = -\frac{2n_1^2 \lambda'}{J} , \quad (5.1.13)$$

where we have set $n_2 = -n_1$ using eqn. (5.1.10). This eigenvalue matches the general

formula in eqn. (5.1.12) restricted to two impurities. The $\mathfrak{su}(2)$ eigenvalue for three impurities with unequal mode indices ($n_1 \neq n_2 \neq n_3$) was calculated in Chapter 4 (and in ref. [38]) and found to be

$$\begin{aligned} \delta E_{S^5}(n_1, n_2, n_3, J) = & -\frac{1}{J\omega_{n_1}\omega_{n_2}\omega_{n_3}} \left\{ [n_1n_2 + n_2^2 + n_1^2(1 + n_2^2\lambda')] \omega_{n_3} \right. \\ & + [n_1n_3 + n_3^2 + n_1^2(1 + n_3^2\lambda')] \omega_{n_2} + [n_2n_3 + n_3^2 + n_2^2(1 + n_3^2\lambda')] \omega_{n_1} \\ & \left. - [n_2n_3 + n_1(n_2 + n_3)] \lambda' \omega_{n_1}\omega_{n_2}\omega_{n_3} \right\} . \end{aligned} \quad (5.1.14)$$

It is also easy to check that eqn. (5.1.12) reproduces this formula exactly for $N_B = 3$.

Since eqn. (5.1.12) matches all previously computed results from the string theory in this sector, it must therefore agree with corresponding $\mathfrak{su}(2)$ gauge theory predictions only to two-loop order in λ . We note, however, that eqn. (5.1.12) is *identical* to the N -impurity $O(J^{-1})$ energy shift (with unequal mode numbers) obtained from the $\mathfrak{su}(2)$ string Bethe ansatz of [44].

To treat the slightly more complicated scenario of overlapping mode indices (which can occur for three or more impurities), we introduce the normalized eigenvectors

$$\frac{1}{\sqrt{N_q!}} (a_q^\dagger)^{N_q} a_{n_1}^\dagger a_{n_2}^\dagger \dots a_{n_{(N_B-N_q)}}^\dagger |J\rangle , \quad (5.1.15)$$

which contain a single subset of N_q bosonic oscillators a_q^\dagger that all share the same mode index q . The remaining indices $n_i \in \{n_1, n_2, \dots, n_{N_B-N_q}\}$ are all separate from q and unequal from each other, such that the level-matching condition in eqn. (5.1.10) now reads

$$N_q q + \sum_{j=1}^{N_B-N_q} n_j = 0 . \quad (5.1.16)$$

For this case we compute a matrix element analogous to that in eqn. (5.1.11):

$$\begin{aligned}
& \frac{1}{N_q!} \langle J | (a_q)^{N_q} a_{n_1} a_{n_2} \dots a_{n_{(N_B-N_q)}} (a_{-n}^\dagger a_{-l}^\dagger a_m a_p) (a_q^\dagger)^{N_q} a_{n_1}^\dagger a_{n_2}^\dagger \dots a_{n_{(N_B-N_q)}}^\dagger | J \rangle \\
&= N_q(N_q - 1) \delta_{p-q} \delta_{m-q} \delta_{n+q} \delta_{l+q} \\
&+ L \sum_{j=1}^{N_B-N_q} \left(\delta_{p-q} \delta_{n+q} \delta_{m-n_j} \delta_{l+n_j} + \delta_{m-q} \delta_{n+q} \delta_{p-n_j} \delta_{l+n_j} \right. \\
&\quad \left. + \delta_{p-q} \delta_{l+q} \delta_{m-n_j} \delta_{n+n_j} + \delta_{m-q} \delta_{l+q} \delta_{p-n_j} \delta_{n+n_j} \right) \\
&+ \frac{1}{2} \sum_{\substack{j,k=1 \\ j \neq k}}^{N_B-N_q} \left(\delta_{n_j+n} \delta_{n_k+l} \delta_{n_j-m} \delta_{n_k-p} + \delta_{n_j+n} \delta_{n_k+l} \delta_{n_k-m} \delta_{n_j-p} \right. \\
&\quad \left. + \delta_{n_j+l} \delta_{n_k+n} \delta_{n_j-m} \delta_{n_k-p} + \delta_{n_j+l} \delta_{n_k+n} \delta_{n_k-m} \delta_{n_j-p} \right) . \tag{5.1.17}
\end{aligned}$$

Using this result, we arrive at the $\mathfrak{su}(2)$ energy shift for string states with N_B total excitations containing an N_q -component subset of oscillators that share the same mode index q :

$$\begin{aligned}
\delta E_{S^5}(\{n_i\}, q, N_q, N_B, J) &= -\frac{N_q(N_q - 1)q^2}{2J\omega_q^2} \\
&- \sum_{j=1}^{N_B-N_q} \frac{N_q}{J\omega_q\omega_{n_j}} [q^2 + n_j^2(1 + q^2\lambda') + q n_j (1 - \omega_q\omega_{n_j}\lambda')] \\
&- \sum_{\substack{j,k=1 \\ j \neq k}}^{N_B-N_q} \frac{1}{2J\omega_j\omega_k} [n_k^2 + n_j^2(1 + n_k^2\lambda') + n_j n_k (1 - \omega_j\omega_k\lambda')] . \tag{5.1.18}
\end{aligned}$$

This formula can be compared with the three-impurity $\mathfrak{su}(2)$ energy shift with two equal mode indices ($N_q = 2$) obtained in Chapter 4. For this particular case we can set the isolated mode number to $-2q$ using the level-matching condition to simplify the result:

$$\delta E_{S^5}(q, J) = -\frac{q^2}{J\omega_q^2\omega_{2q}} [\omega_{2q} (5 + 4q^2\lambda') + \omega_q (6 + 8q^2\lambda')] . \tag{5.1.19}$$

It is easy to show that eqn. (5.1.18) exactly reproduces this energy shift when re-

stricted to $N_B = 3$ with a subset of two mode numbers equal to q .

We now generalize the analysis completely by using eigenstates with M mode-index subsets, where all mode indices are equal within these subsets:

$$\frac{(a_{q_1}^\dagger)^{N_{q_1}}}{\sqrt{N_{q_1}!}} \frac{(a_{q_2}^\dagger)^{N_{q_2}}}{\sqrt{N_{q_2}!}} \cdots \frac{(a_{q_M}^\dagger)^{N_{q_M}}}{\sqrt{N_{q_M}!}} |J\rangle .$$

The j^{th} subset contains N_{q_j} oscillators with equal mode index q_j , and the total impurity number is again N_B , such that

$$\sum_{i=1}^M N_{q_i} = N_B , \quad \sum_{i=1}^M N_{q_i} q_i = 0 . \quad (5.1.20)$$

The matrix element of $a_{-n}^\dagger a_{-l}^\dagger a_m a_p$ between the above states, analogous to eqns. (5.1.11, 5.1.17), is

$$\begin{aligned} \langle J | & \frac{(a_{q_1})^{N_{q_1}}}{\sqrt{N_{q_1}!}} \cdots \frac{(a_{q_M})^{N_{q_M}}}{\sqrt{N_{q_M}!}} \left(a_{-n}^\dagger a_{-l}^\dagger a_m a_p \right) \frac{(a_{q_1}^\dagger)^{N_{q_1}}}{\sqrt{N_{q_1}!}} \cdots \frac{(a_{q_M}^\dagger)^{N_{q_M}}}{\sqrt{N_{q_M}!}} |J\rangle \\ &= \sum_{j=1}^M N_{q_j} (N_{q_j} - 1) \delta_{n+n_j} \delta_{l+n_j} \delta_{m-n_j} \delta_{p-n_j} + \frac{1}{2} \sum_{\substack{j,k=1 \\ j \neq k}}^M N_{q_j} N_{q_k} \left(\delta_{n+n_k} \delta_{l+n_j} \delta_{m-n_k} \delta_{p-n_j} \right. \\ & \quad \left. + \delta_{n+n_j} \delta_{l+n_k} \delta_{m-n_k} \delta_{p-n_j} + \delta_{n+n_k} \delta_{l+n_j} \delta_{m-n_j} \delta_{p-n_k} + \delta_{n+n_j} \delta_{l+n_k} \delta_{m-n_j} \delta_{p-n_k} \right) . \end{aligned} \quad (5.1.21)$$

We thereby obtain the completely general $\mathfrak{su}(2)$ energy shift for N_B -impurity string states containing M equal-mode-index subsets of oscillators:

$$\begin{aligned} \delta E_{S^5}(\{q_i\}, \{N_{q_i}\}, M, J) &= -\frac{1}{2J} \left\{ \sum_{j=1}^M N_{q_j} (N_{q_j} - 1) \left(1 - \frac{1}{\omega_{q_j}^2 \lambda'} \right) \right. \\ & \quad \left. + \sum_{\substack{j,k=1 \\ j \neq k}}^M \frac{N_{q_j} N_{q_k}}{\omega_{q_j} \omega_{q_k}} [q_k^2 + q_j^2 \omega_{q_k}^2 \lambda' + q_j q_k (1 - \omega_{q_j} \omega_{q_k} \lambda')] \right\} . \end{aligned} \quad (5.1.22)$$

This master formula can be used to determine the $\mathfrak{su}(2)$ string energy spectrum to $O(J^{-1})$ for all possible physical string states in this sector.

By taking $M = 2$ and setting $N_{n_1} = N_{n_2} = 1$ (using the unequal mode indices $\{n_1, n_2\}$), we recover from this equation the exact two-impurity result recorded in eqn. (5.1.13) above, with $n_2 = -n_1$. For $M = 3$ and $N_{n_1} = N_{n_2} = N_{n_3} = 1$, we get the complete three-impurity unequal-mode-number ($n_1 \neq n_2 \neq n_3$) formula found in eqn. (5.1.14). Finally, the three-impurity eigenvalue with two equal mode indices ($q_1 = q_2$, $q_3 = -2q_1$) given in eqn. (5.1.19) can also be extracted from eqn. (5.1.22) by setting $M = 2$, $N_{q_1} = 2$ and $N_{q_2} = 1$.

We also note that eqn. (5.1.22) agrees perfectly with the corresponding near-pp-wave formula derived from the $\mathfrak{su}(2)$ string Bethe ansatz of [44] for completely general mode-number assignment. This successful match stands as very strong evidence that their ansatz is correct, at least to $O(J^{-1})$.

5.1.2 The $SO(4)_{AdS}$ ($\mathfrak{sl}(2)$) sector

Following the derivation of eqn. (5.1.22) for the energy eigenvalues of arbitrary string states in the symmetric-traceless $SO(4)_{S^5}$ sector, it is straightforward to find the analogous expression for symmetric-traceless string states excited in the $SO(4)_{AdS}$ subspace, dual to operators in the $\mathfrak{sl}(2)$ sector of the corresponding gauge theory. We can define, for example,

$$a_n = \frac{1}{\sqrt{2}} (a_n^1 + i a_n^2) \ , \quad \bar{a}_n = \frac{1}{\sqrt{2}} (a_n^1 - i a_n^2) \ , \quad (5.1.23)$$

and carry out the above calculations by computing general matrix elements of $a_{-n}^\dagger a_{-l}^\dagger a_m a_p$ defined in terms of these oscillators. (Here we can project onto any (n, m) -plane in the AdS_5 subspace, as long as $n \neq m$.) General string energy eigenvalues in the

$SO(4)_{AdS}$ symmetric-traceless irrep are thus found to be

$$\begin{aligned} \delta E_{AdS}(\{q_i\}, \{N_{q_i}\}, M, J) = & \frac{1}{2J} \left\{ \sum_{j=1}^M N_{q_j} (N_{q_j} - 1) \left(1 - \frac{1}{\omega_{q_j}^2 \lambda'} \right) \right. \\ & \left. + \sum_{\substack{j,k=1 \\ j \neq k}}^M \frac{N_{q_j} N_{q_k}}{\omega_{q_j} \omega_{q_k}} q_j q_k [1 - q_j q_k \lambda' + \omega_{q_j} \omega_{q_k} \lambda'] \right\}. \end{aligned} \quad (5.1.24)$$

For later reference we record the limit of this equation for states with completely unequal mode indices ($\{N_{n_i}\} = 1, M = N_B$):

$$\delta E_{AdS}(\{n_i\}, N_B, J) = \frac{1}{2J} \sum_{\substack{j,k=1 \\ j \neq k}}^{N_B} \frac{n_j n_k}{\omega_{n_j} \omega_{n_k}} [1 - n_j n_k \lambda' + \omega_{n_j} \omega_{n_k} \lambda'] . \quad (5.1.25)$$

When $M = 2$ and $N_{n_1} = N_{n_2} = 1$ in eqn. (5.1.25), we find the two-impurity eigenvalue (with $n_2 = -n_1$)

$$\delta E_{AdS}(n_1, J) = -\frac{2 n_1^2 \lambda'}{J}, \quad (5.1.26)$$

which agrees with the two-impurity result reported in Chapter 3 [26,27] (the $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ eigenvalues are degenerate in the two-impurity regime). For the three-impurity eigenvalue with three unequal mode indices we set $M = 3$ and $N_{n_1} = N_{n_2} = N_{n_3} = 1$ to obtain

$$\begin{aligned} \delta E_{AdS}(n_1, n_2, n_3, J) = & \frac{1}{J \omega_{n_1} \omega_{n_2} \omega_{n_3}} \left\{ n_1 n_3 (1 - n_1 n_3 \lambda') \omega_{n_2} + n_1 n_2 (1 - n_1 n_2 \lambda') \omega_{n_3} \right. \\ & \left. + n_2 n_3 (1 - n_2 n_3 \lambda') \omega_{n_1} + [n_1 n_2 + n_3 (n_1 + n_2)] \lambda' \omega_{n_1} \omega_{n_2} \omega_{n_3} \right\}, \end{aligned} \quad (5.1.27)$$

which precisely reproduces the corresponding $\mathfrak{sl}(2)$ result reported in Chapter 4 [38]. Finally, by setting $M = 2, N_{q_1} = 2, N_{q_2} = 1$ and $q_1 = q_2 = q, q_3 = -2q$, eqn. (5.1.24)

provides the following three-impurity eigenvalue with two equal mode indices:

$$\delta E_{AdS}(q, J) = -\frac{q^2}{J\omega_q^2\omega_{2q}} \left[\omega_{2q} (3 + 4q^2\lambda') + \omega_q (4 + 8q^2\lambda') \right] . \quad (5.1.28)$$

This again matches the corresponding three-impurity formula computed in Chapter 4.

5.1.3 The $\mathfrak{su}(1|1)$ sector

Based on the above results in the bosonic $SO(4)_{AdS}$ and $SO(4)_{S^5}$ symmetric-traceless sectors, we can easily formulate a conjecture for the N -impurity eigenvalue of symmetrized pure-fermion states in either the $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ of $SO(4) \times SO(4)$, labeled by the $\mathfrak{su}(1|1)$ subalgebra. We first note that, since these states are composed of fermionic oscillators that are symmetrized in their spinor indices, no states in this sector can carry subsets of overlapping mode numbers (since they would automatically vanish). Furthermore, when restricting to states with completely unequal mode indices, we can see that the N -impurity eigenvalues obtained for the $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ sectors (eqns. (5.1.12) and (5.1.25)) are obvious generalizations of the corresponding three-impurity formulas (eqns. (5.1.14) and (5.1.27), respectively). Namely, if the three-impurity eigenvalues take the generic form

$$\delta E(n_1, n_2, n_3, J) = \sum_{\substack{j,k=1 \\ j \neq k}}^3 F(n_j, n_k) , \quad (5.1.29)$$

the N -impurity generalization is simply

$$\delta E(\{n_i\}, N, J) = \sum_{\substack{j,k=1 \\ j \neq k}}^N F(n_j, n_k) . \quad (5.1.30)$$

By carrying this over to the $\mathfrak{su}(1|1)$ sector, we find the N -impurity eigenvalue of H_{FF} between symmetrized $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ fermions (the eigenvalues of both are

necessarily degenerate):

$$\delta E_{\mathfrak{su}(1|1)}(\{n_i\}, N_F, J) = -\frac{1}{4J} \sum_{\substack{j,k=1 \\ j \neq k}}^{N_F} \frac{1}{\omega_{n_j} \omega_{n_k}} \left[n_j^2 + n_k^2 + 2n_j^2 n_k^2 \lambda' - 2n_j n_k \omega_{n_j} \omega_{n_k} \lambda' \right] . \quad (5.1.31)$$

For $N_F = 2$, this formula matches the two-impurity result in Chapter 3:

$$\delta E_{\mathfrak{su}(1|1)}(n_1, J) = -\frac{2n_1^2 \lambda'}{J} , \quad (5.1.32)$$

with $n_2 = -n_1$ (this eigenvalue overlaps with the corresponding two-impurity $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ values). When $N_F = 3$ we of course recover the three-impurity eigenvalue reported in Chapter 4:

$$\begin{aligned} \delta E_{\mathfrak{su}(1|1)}(n_1, n_2, n_3, J) = & -\frac{1}{4J\omega_{n_1}\omega_{n_2}\omega_{n_3}} \left\{ -4(n_2 n_3 + n_1(n_2 + n_3)) \lambda' \omega_{n_1} \omega_{n_2} \omega_{n_3} \right. \\ & \left. + \left[\omega_{n_1} (2n_3^2 + 4n_2^2 n_3^2 \lambda' + 2n_2^2) + (n_3 \rightarrow n_2, n_2 \rightarrow n_1, n_1 \rightarrow n_3) + (n_1 \rightleftharpoons n_2) \right] \right\} . \end{aligned} \quad (5.1.33)$$

It would be straightforward to check eqn. (5.1.31) against an explicit four-impurity calculation in the string theory, for example. Better yet, one might carry out the direct N -impurity calculation in the H_{FF} sector analogous to the above calculations for H_{BB} . The latter would certainly be more technically complicated than in the bosonic sectors, and for the moment we leave eqn. (5.1.31) as it stands, withholding direct verification for a future study.

5.2 Spectral decomposition

At one- and two-loop order in λ' we can infer from basic arguments the spectral decomposition of the extended N -impurity superconformal multiplet of $O(J^{-1})$ energy corrections to the pp-wave limit. For simplicity we will restrict the discussion to

eigensystems with completely unequal mode numbers, though the generalization to more complicated cases is straightforward. To begin we will review the two- and three-impurity supermultiplet structures studied in [26, 27, 38].

We denote the one- and two-loop energy eigenvalue shifts as $\Lambda^{(1)}$ and $\Lambda^{(2)}$, according to the generic formula

$$\begin{aligned} E(\{n_j\}, N, J) = & N + \frac{\lambda'}{2} \sum_{j=1}^N n_j^2 \left(1 + \frac{\Lambda^{(1)}}{J} + O(J^{-2}) \right) \\ & - \frac{\lambda'^2}{4} \sum_{j=1}^N n_j^4 \left(\frac{1}{2} + \frac{\Lambda^{(2)}}{J} + O(J^{-2}) \right) + O(\lambda'^3) . \end{aligned} \quad (5.2.1)$$

The fact that these energy shifts can be expressed as coefficients of $\sum n_j^2$ and $\sum n_j^4$ is not obvious. In the two- and three-impurity cases this was shown to be true by direct diagonalization of the Hamiltonian. By expanding eqns. (5.1.22, 5.1.24, 5.1.31) in small λ' , it can also be seen that the more general N -impurity $\mathfrak{su}(2)$, $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ eigenvalues adhere to this structure to two-loop order. We will argue that the remaining energy shifts (those in non-protected subsectors) can be obtained from the protected sectors through half-integer shifts of the S^5 angular momentum J : it will therefore be seen that all energies considered here will appear in the form given in eqn. (5.2.1).

As described above, the conformal invariance of the full $\mathfrak{psu}(2, 2|4)$ symmetry algebra of the theory guarantees that the energy eigenvalues (and hence $\Lambda^{(1)}$ and $\Lambda^{(2)}$) will be organized into conformal (sub)multiplets built on conformal primary (or highest weight) states. For the sake of continuity we will briefly review this here. Within a given submultiplet we refer to states with lowest energy as super-primary states, and the other conformal primaries within the submultiplet are obtained by acting on super-primaries with any of the eight supercharges, labeled by \mathcal{Q}_α , that increment $\Lambda^{(1)}$ or $\Lambda^{(2)}$ by a fixed amount but leave the impurity number unchanged. In the gauge theory these charges are understood to shift both the operator dimension and R -charge such that $\Delta = D - R$ remains fixed within the submultiplet. Acting

with L_{sub} factors of these supercharges on a super-primary generates nine levels within each submultiplet labeled by $L_{\text{sub}} = 0, \dots, 8$. If the lowest energy level ($L_{\text{sub}} = 0$) in the submultiplet is occupied by p degenerate super-primaries, the $L_{\text{sub}}^{\text{th}}$ level will therefore contain $p C_{L_{\text{sub}}}^8$ degenerate states, where C_n^m is the binomial coefficient. Furthermore, if the super-primary in a given submultiplet is a spacetime boson, the $L_{\text{sub}} = \text{even}$ levels of the submultiplet will all be bosonic, and the $L_{\text{sub}} = \text{odd}$ levels will be fermionic. The opposite is true if the bottom state is fermionic.

As an example, consider the one-loop, two-impurity supermultiplet structure studied in Chapter 3. The spectrum in this case contains only a single multiplet built on a scalar super-primary (labeled by 1_B , where the subscript denotes a bosonic level) with $O(1/J)$ one-loop energy shift $\Lambda^{(1)} = -6$. The $L_{\text{sub}} = 1$ level therefore has eight degenerate states (8_F) with $\Lambda^{(1)} = -5$, the $L_{\text{sub}} = 2$ level contains 28_B states with $\Lambda^{(1)} = -4$ and so on. We record the two-impurity supermultiplet structure in table 5.2 for comparison with higher-impurity spectra. The one-loop energies of the

L_{sub}	0	1	2	3	4	5	6	7	8
	1_B	8_F	28_B	56_F	70_B	56_F	28_B	8_F	1_B
$\Lambda^{(1)}(L_{\text{sub}})$	-6	-5	-4	-3	-2	-1	0	1	2
$\Lambda^{(2)}(L_{\text{sub}})$	-4	-3	-2	-1	0	1	2	3	4

Table 5.2: Submultiplet breakup of the 256-dimensional two-impurity spectrum

three protected $\mathfrak{sl}(2)$, $\mathfrak{su}(2)$ and $\mathfrak{su}(1|1)$ subsectors studied here are degenerate in the two-impurity regime and lie in the boxed 70_B “centroid” level in table 5.2. We also record in table 5.2 the two-loop energy shifts $\Lambda^{(2)}$, which are offset from the one-loop values by two: $\Lambda^{(2)} = \Lambda^{(1)} + 2$.

In the gauge theory there are 16 operators that increment the impurity number by one and shift the R -charge by certain amounts [28]. Four of these act on single-trace operators by rotating the $SO(6)$ scalars Z (carrying one unit of R -charge) into ϕ (which carry zero R -charge): they increase the operator impurity number by one and decrease the R -charge by one ($N \rightarrow N + 1$, $R \rightarrow R - 1$). Four operators rotate Z

into ∇Z , increasing N by one and leaving the R -charge fixed. The remaining eight operators are fermionic and send $N \rightarrow N+1$, $R \rightarrow R+1/2$. If one uses these operators to generate N -impurity super-primaries from those in the $(N-1)$ -impurity spectrum, an immediate implication is that, within a given N -impurity spectrum of anomalous dimensions, all of the eigenvalues in the gauge theory will be related to each other by half-integer shifts in the R -charge. Certain energy levels will therefore be common to all of the submultiplets in the spectrum built on super-primary operators, and this special degeneracy can be used to deduce the overall structure of the extended supermultiplet. This degeneracy, however, only persists in the string theory to two-loop order in λ' , and it is for this reason that we are forced to limit the general superstring spectral decomposition to two-loop order in the expansion. (It will be shown below, however, that a certain subset of submultiplets in the string theory can always be determined to *all* orders in λ' .)

Sending $J \rightarrow J + A$ on the string side (dual to an R -charge shift in the gauge theory) shifts $\Lambda^{(1)}$ and $\Lambda^{(2)}$ by $-2A$: starting from the two-impurity super-primary (1_B) with energy $\Lambda^{(1)} = -6$, the string versions of the 16 impurity-increasing operators can be understood to generate four (degenerate) bosonic three-impurity super-primaries with $\Lambda^{(1)} = -8$, eight fermionic three-impurity super-primaries with $\Lambda^{(1)} = -7$ and four bosonic three-impurity super-primaries with $\Lambda^{(1)} = -6$. By acting with the eight charges \mathcal{Q}_α we then generate submultiplets based on each of these super-primaries whose levels are populated by $p C_{L_{\text{sub}}}^8$ degenerate states, where p here is either four (for the two four-dimensional bosonic super-primary levels) or eight (for the eight-dimensional fermionic super-primary level). The submultiplets themselves can be labeled by a separate index L' , in this case running over $L' = 0, \dots, 2$.

The complete three-impurity multiplet structure is recorded in table 5.3. Here there are a total of 11 levels in the extended supermultiplet, and we label these with the index L such that $L = L_{\text{sub}} + L'$. In table 5.3 the closed $\mathfrak{su}(2)$ sector lies in the boxed 280_B level in the $L' = 0$ submultiplet with $\Lambda^{(1)} = -4$, the $\mathfrak{sl}(2)$ eigenvalue ($\Lambda^{(1)} = -2$) is in the boxed 280_B level of the $L' = 2$ submultiplet and

the $\mathfrak{su}(1|1)$ eigenvalue ($\Lambda^{(1)} = -3$) is in the 560_F level of the $L' = 1$ submultiplet. For any impurity number these protected eigenvalues will always lie at the $L_{\text{sub}} = 4$ level within their respective submultiplets. We also note that, in the L' direction, the $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ eigenvalues will correspond to eigenstates composed purely of S^5 or AdS_5 bosonic excitations, and will therefore fall into the “bottom” and “top” submultiplets, respectively (the $L' = 0$ and $L' = 2$ levels in the three-impurity case). Similarly, the $\mathfrak{su}(1|1)$ eigenvalue will correspond to eigenstates composed of either $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ excitations, and always lie in the “centroid” submultiplet in the L' direction (the $L' = 1$ level for three impurities). The energies shared by each of the submultiplets can be collected into degenerate levels of the complete supermultiplet. This total level degeneracy $D(L)$ is recorded in the bottom row of table 5.3.

$L' \backslash L$	0	1	2	3	4	5	6	7	8	9	10
0	4	32	112	224	280	224	112	32	4		
1		8	64	224	448	560	448	224	64	8	
2			4	32	112	224	280	224	112	32	4
$\Lambda^{(1)}(L)$	-8	-7	-6	-5	-4	-3	-2	-1	0	1	2
$\Lambda^{(2)}(L)$	-6	-5	-4	-3	-2	-1	0	1	2	3	4
$D(L)$	4_B	40_F	180_B	480_F	840_B	1008_F	840_B	480_F	180_B	40_F	4_B

Table 5.3: Submultiplet breakup of the 4,096-dimensional three-impurity spectrum

It is easy to generalize this supermultiplet structure to arbitrary impurity number based on how the complete three-impurity spectrum is generated from the two-impurity supermultiplet above. For N impurities, the complete supermultiplet will have a total of 16^N states and $5+2N$ levels: the supermultiplet level index L therefore runs over $L = 0, \dots, (4+2N)$. The entire supermultiplet breaks into $2N-3$ submultiplets, each of which have nine sub-levels labeled by $L_{\text{sub}} = 0, \dots, 8$. The submultiplets themselves are labeled by the index L' , which runs over $L' = 0, \dots, (2N-4)$. The

one-loop energy shifts within the L'^{th} submultiplet at level L_{sub} are thus given by

$$\Lambda_{\text{sub}}^{(1)}(L', L_{\text{sub}}, N) = L' + L_{\text{sub}} - 2(N + 1) . \quad (5.2.2)$$

Equivalently, the L^{th} level of the entire supermultiplet has energy shift

$$\Lambda^{(1)}(L, N) = L - 2(N + 1) . \quad (5.2.3)$$

The number of degenerate states at level L_{sub} within the L'^{th} submultiplet is

$$D_{\text{sub}}(L', L_{\text{sub}}, N) = 4^{N-2} C_{L'}^{2N-4} C_{L_{\text{sub}}}^8 , \quad (5.2.4)$$

so that the total dimension of the L'^{th} submultiplet is $256 \times 4^{N-2} C_{L'}^{2N-4}$. By summing the submultiplet degeneracies over a given supermultiplet level L , the total number of degenerate states at level L in the supermultiplet is given (in terms of Euler's Γ function) by

$$D(L, N) = \frac{4^{N-2} \Gamma(2N + 5)}{\Gamma(2N + 5 - L) \Gamma(1 + L)} . \quad (5.2.5)$$

The level is bosonic when L is even and fermionic when L is odd. As a verification of this formula, we can check that the total number of states in the N -impurity supermultiplet is indeed

$$\sum_{L=0}^{4+2N} \frac{4^{N-2} \Gamma(2N + 5)}{\Gamma(2N + 5 - L) \Gamma(1 + L)} = 16^N . \quad (5.2.6)$$

As noted above, the one-loop N -impurity $\mathfrak{su}(2)$ energy corresponds to eigenstates that are composed purely of symmetric-traceless $(\mathbf{1}, \mathbf{1}; \mathbf{2}, \mathbf{2})$ excitations: since each of these excitations increments the angular momentum J by one, the energy eigenvalue must therefore lie within a submultiplet built on super-primary states that exhibit the lowest possible energy in the extended supermultiplet. In other words, the $\mathfrak{su}(2)$

eigenvalue always lies at level $L_{\text{sub}} = 4$ of the $L' = 0$ submultiplet and, using the general formula in eqn. (5.2.2), we see that it exhibits the one-loop energy shift

$$\Lambda_{S^5}^{(1)}(N) = \Lambda_{\text{sub}}^{(1)}(L' = 0, L_{\text{sub}} = 4, N) = -2(N - 1) . \quad (5.2.7)$$

As a cross-check on this result, we note that this agrees with the one-loop limit of the general $\mathfrak{su}(2)$ eigenvalue formula (with unequal mode indices) in eqn. (5.1.12) above (with $N_B = N$):

$$\delta E_{S^5}(\{n_i\}, N, J) = -\frac{1}{2J} \sum_{\substack{j,k=1 \\ j \neq k}}^N (n_j^2 + n_k^2) \lambda' + O(\lambda'^2) = -\frac{1}{J} \sum_{j=1}^N (N-1) n_j^2 \lambda' + O(\lambda'^2) . \quad (5.2.8)$$

(Note the prefactor of $1/2$ in the definition of $\Lambda^{(1)}$ in eqn. (5.2.1).) At this point we also see that $\Lambda^{(1)}$ indeed appears as a coefficient of $\sum n_j^2$, as given in eqn. (5.2.1).

The N -impurity $\mathfrak{sl}(2)$ eigenvalue, composed entirely of $(\mathbf{2}, \mathbf{2}; \mathbf{1}, \mathbf{1})$ excitations, must lie in the “top” $L' = 2N - 4$ submultiplet at $L_{\text{sub}} = 4$. This gives the one-loop energy shift

$$\Lambda_{AdS}^{(1)}(N) = -2 . \quad (5.2.9)$$

To check this we use the general $\mathfrak{sl}(2)$ formula for completely unequal mode indices in eqn. (5.1.25), and again expand to one-loop order in λ' :

$$\delta E_{AdS}(\{n_i\}, N, J) = \frac{1}{J} \sum_{\substack{j,k=1 \\ j \neq k}}^N n_j n_k \lambda' + O(\lambda'^2) . \quad (5.2.10)$$

With the level-matching condition $\sum_{j=1}^N n_j = 0$ this becomes

$$\delta E_{AdS}(\{n_i\}, N, J) = -\frac{1}{J} \sum_{j=1}^N n_j^2 \lambda' + O(\lambda'^2) , \quad (5.2.11)$$

which agrees perfectly with the prediction in eqn. (5.2.9) (and again confirms that $\Lambda^{(1)}$ here is a coefficient of $\sum n_j^2$).

Finally, the $\mathfrak{su}(1|1)$ one-loop eigenvalue, composed of either $(\mathbf{2}, \mathbf{1}; \mathbf{2}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{2}; \mathbf{1}, \mathbf{2})$ spinors, lies in the $L' = N - 2$ submultiplet at $L_{\text{sub}} = 4$, exhibiting the one-loop energy shift

$$\Lambda_{\mathfrak{su}(1|1)}^{(1)}(N) = -N . \quad (5.2.12)$$

Using eqn. (5.1.31) we see that

$$\begin{aligned} \delta E_{\mathfrak{su}(1|1)}(\{n_i\}, N, J) &= -\frac{1}{4J} \sum_{\substack{j,k=1 \\ j \neq k}}^N (n_j - n_k)^2 \lambda' + O(\lambda'^2) \\ &= -\frac{1}{2J} \sum_{j=1}^N N n_j^2 \lambda' , \end{aligned} \quad (5.2.13)$$

where we have again invoked the level-matching condition to derive the last line. This of course agrees with eqn. (5.2.12). For reference we present in table 6.4 the complete 65,536-dimensional four-impurity spectrum of one- and two-loop energies. The $\mathfrak{su}(2)$ eigenvalue in this case lies in the boxed 1120_B level with $\Lambda^{(1)} = -6$, the $\mathfrak{su}(1|1)$ eigenvalue is in the 6720_B level with $\Lambda^{(1)} = -4$, and the $\mathfrak{sl}(2)$ energy lies in the 1120_B level with $\Lambda^{(1)} = -2$.

$L' \setminus L$	0	1	2
0	16	128	448
1		64	512
2			96
3			
4			
$\Lambda^{(1)}(L)$	-10	-9	-8
$\Lambda^{(2)}(L)$	-8	-7	-6
$D(L)$	16_B	192_F	1056_B

Table 5.4:

Comparing the $\Lambda^{(2)}$ and $\Lambda^{(1)}$ spectra in tables 5.2 and 5.3 (which are determined directly from the string Hamiltonian), we see that the spectrum of $\Lambda^{(2)}$ is identical to $\Lambda^{(1)}$ up to an overall shift. The two-loop analogue of the general N -impurity energy shift of eqn. (5.2.2) is therefore

$$\Lambda_{\text{sub}}^{(2)}(L', L_{\text{sub}}, N) = L' + L_{\text{sub}} - 2N . \quad (5.2.14)$$

Equivalently, we have $\Lambda^{(2)}(L, N) = L - 2N$ for the entire supermultiplet shift in terms of L .

Similar to the one-loop case, we can test this two-loop formula using the N -impurity results derived above in the three protected sectors. According to eqn. (5.2.14), the $\mathfrak{su}(2)$ eigenvalue in the $L' = 0$ submultiplet at level $L_{\text{sub}} = 4$ has the following two-loop energy shift:

$$\Lambda_{S^5}^{(2)}(N) = 4 - 2N . \quad (5.2.15)$$

Isolating the two-loop energy eigenvalue $\delta E_{S^5}^{(2)}$ from the N -impurity $\mathfrak{su}(2)$ equation (5.1.12), we have

$$\begin{aligned} \delta E_{S^5}^{(2)}(\{n_i\}, N, J) &= \frac{1}{4J} \sum_{\substack{j,k=1 \\ j \neq k}}^N (n_j^4 + n_j^3 n_k + n_j n_k^3 + n_k^4) \lambda'^2 \\ &= -\frac{1}{4J} \sum_{j=1}^N (n_j^4) (4 - 2N) \lambda'^2 , \end{aligned} \quad (5.2.16)$$

which matches our prediction. The $\mathfrak{sl}(2)$ eigenvalue in the $L' = 2N - 4$ submultiplet is predicted to vanish

$$\Lambda_{AdS}^{(2)}(N) = 0 , \quad (5.2.17)$$

which agrees with the two-loop expansion term in eqn. (5.1.25):

$$\delta E_{AdS}^{(2)}(\{n_i\}, J) = -\frac{1}{4J} \sum_{\substack{j,k=1 \\ j \neq k}}^N [n_j n_k (n_j + n_k)^2] \lambda'^2 = 0 . \quad (5.2.18)$$

Finally, the $\mathfrak{su}(1|1)$ pure-fermion sector in the $L' = N - 2$ submultiplet at $L_{\text{sub}} = 4$ should have an energy shift of

$$\Lambda_{\mathfrak{su}(1|1)}^{(2)}(N) = 2 - N , \quad (5.2.19)$$

which agrees with the $\mathfrak{su}(1|1)$ formula given in eqn. (5.1.31):

$$\begin{aligned} \delta E_{\mathfrak{su}(1|1)}^{(2)}(\{n_i\}, N, J) &= \frac{1}{8J} \sum_{\substack{j,k=1 \\ j \neq k}}^N (n_j^2 - n_k^2)^2 \lambda'^2 \\ &= -\frac{1}{4J} \sum_{j=1}^N (n_j^4) (2 - N) \lambda'^2 . \end{aligned} \quad (5.2.20)$$

As described in Chapter 4, it should also be noted that since we know the $\mathfrak{su}(2)$, $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ eigenvalues to all orders in λ' , we can easily determine complete all-loop energy formulas for the three submultiplets to which these eigenvalues belong. It was previously noted that the eight supercharges (\mathcal{Q}_α) that act as raising operators within each submultiplet are known in the gauge theory to shift both the dimension and R -charge by $1/2$ such that $\Delta = D - R$ is kept fixed. Because all states within a given submultiplet share the same Δ , the string energy shift at any level L_{sub} can therefore be obtained from that at some level L'_{sub} (not to be confused with L') by replacing

$$J \rightarrow J - L_{\text{sub}}/2 + L'_{\text{sub}}/2$$

in the energy eigenvalue evaluated at sub-level L'_{sub} . Since we are expanding to $O(J^{-1})$, however, this replacement can only affect the eigenvalues δE via the $O(J^0)$

BMN term in the pp-wave limit. For the protected eigenvalues determined above at $L_{\text{sub}} = 4$, we therefore find the all-loop energy shift for the entire submultiplet by including the appropriate $O(J^{-1})$ contribution from the BMN formula

$$E_{\text{BMN}} = \sum_{j=1}^N \sqrt{1 + \frac{n_j^2 \lambda}{(J + 2 - L_{\text{sub}}/2)^2}} . \quad (5.2.21)$$

Explicitly, the complete level spectra of the $L' = 0$, $L' = N - 2$ and $L' = 2N - 4$ submultiplets are given, to all orders in λ' , by

$$\delta E(\{n_j\}, L_{\text{sub}}, N, J) = \frac{\lambda'}{2J} \sum_{j=1}^N \frac{n_j^2 (L_{\text{sub}} - 4)}{\sqrt{1 + n_j^2 \lambda'}} + \delta E_{L_{\text{sub}}=4}(\{n_j\}, J) , \quad (5.2.22)$$

where $\delta E_{L_{\text{sub}}=4}$ is the $L_{\text{sub}} = 4$ energy shift in the submultiplet of interest. Since the level degeneracy among submultiplets is generally broken beyond two-loop order, it is difficult to obtain similar expressions for submultiplets not containing the $\mathfrak{su}(2)$, $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ protected eigenvalues. This can possibly be addressed by relying directly on the commutator algebra of various impurity-increasing operators in the string theory, and we will return to this problem in a future study.

5.3 Discussion

In this chapter we have directly computed the near-pp-wave eigenvalues of N -impurity bosonic string states with arbitrary mode-number assignment lying in the protected symmetric-traceless irreps of the AdS_5 ($\mathfrak{sl}(2)$) and S^5 ($\mathfrak{su}(2)$) subspaces. Based on the observation that the $\mathfrak{su}(2)$ and $\mathfrak{sl}(2)$ eigenvalues are simple generalizations of the three-impurity results obtained in Chapter 4, we have also presented a conjecture for the N -impurity eigenvalues of symmetrized-fermion states in the $\mathfrak{su}(1|1)$ sector. This conjecture meets several basic expectations and we believe that it is correct. (It would be satisfying, however, to derive the $\mathfrak{su}(1|1)$ eigenvalue formula directly from the fermionic sector of the string theory.) We have also found that the $\mathfrak{su}(2)$ eigenvalues

perfectly match, to all orders in λ' , the corresponding eigenvalue predictions given by the string Bethe ansatz of [44]. Along these lines, it would be very interesting to have long-range Bethe ansätze analogous to [44] for the entire $\mathfrak{psu}(2, 2|4)$ algebra of the theory.

The supermultiplet decomposition given in Section 5.2 is based on the breakup of the energy spectrum observed between the two- and three-impurity regime, and is precisely what is expected from the gauge theory based on how 16 particular charges are known to act on operators that are dual to the string states of interest [28, 38]. Assuming that this mechanism is not specific to the three-impurity case, we were able to generalize the decomposition of the N -impurity (unequal mode index) supermultiplet to two-loop order in λ' . By knowing where the eigenvalues of the $\mathfrak{su}(2)$, $\mathfrak{sl}(2)$ and $\mathfrak{su}(1|1)$ sectors are supposed to appear in this decomposition, we were able to provide a stringent cross-check of our results, and we have found perfect agreement. Given the many implicit assumptions in this procedure, however, it would be instructive to perform a direct diagonalization of the four-impurity Hamiltonian to test our predictions. While such a test is likely to be computationally intensive, the problem could be simplified to some extent by restricting to the pure-boson H_{BB} sector at one loop in λ' . We of course expect complete agreement with the results presented in this chapter.

Chapter 6

Integrability in the quantum string theory

The emergence of integrable structures from planar $\mathcal{N} = 4$ super-Yang-Mills (SYM) theory and type IIB string theory on $AdS_5 \times S^5$ has renewed hope that 't Hooft's formulation of large- N_c QCD may eventually lead to an exact solution. If both the gauge and string theories are in fact integrable, each will admit infinite towers of hidden charges and, analogous to the usual identification of the string theory Hamiltonian with the gauge theory dilatation generator, there will be an infinite number of mappings between the higher hidden charges of both theories. This has led to many novel tests of the AdS/CFT correspondence, particularly in the context of the pp-wave/BMN limits [22, 25, 97]. Barring an explicit solution, one would hope that both theories will at least be shown to admit identical Bethe ansatz equations, allowing us to explore a much larger region of the gauge/string duality.

As described above, the fact that the gauge theory harbors integrable structures was realized by Minahan and Zarembo when they discovered that a particular $SO(6)$ -invariant sector of the SYM dilatation generator can be mapped, at one-loop order in the 't Hooft coupling $\lambda = g_{\text{YM}}^2 N_c$, to the Hamiltonian of an integrable quantum spin chain with $SO(6)$ vector lattice sites [39]. The Hamiltonian of this system can be diagonalized by solving a set of algebraic Bethe ansatz equations: the problem of computing operator anomalous dimensions in this sector of the gauge theory was thus reduced in [39] to solving the set of Bethe equations specific to the $\mathfrak{so}(6)$ sec-

tor of the theory. The correspondence between operator dimensions and integrable spin chain systems at one loop in λ was extended to include the complete $\mathfrak{psu}(2, 2|4)$ superconformal symmetry algebra of planar $\mathcal{N} = 4$ SYM theory by Beisert and Staudacher in [32]. Studies of higher-loop integrability in the gauge theory were advanced in [34, 37], where so-called long-range Bethe ansatz equations, which are understood to encode interactions on the spin lattice that extend beyond nearest-neighbor sites, were developed for a closed bosonic $\mathfrak{su}(2)$ sector of the gauge theory. The dynamics of the gauge theory therefore appear to be consistent with the expectations of integrability, at least to three-loop order in the 't Hooft expansion, and there is convincing evidence that this extends to even higher order [34, 35].

Concurrent with the introduction of the Bethe ansatz formalism in the $\mathfrak{so}(6)$ sector of the gauge theory [39], related developments emerged from studies of semiclassical configurations of rotating string on $AdS_5 \times S^5$. This branch of investigation began with [98], where the pp-wave limit of the string theory was reinterpreted in the context of a semiclassical expansion about certain solitonic solutions in the full $AdS_5 \times S^5$ target space. Using this semiclassical picture, Frolov and Tseytlin computed a class of two-spin string solutions in [99], demonstrating explicitly how stringy corrections in the large-spin limit give rise to systems that can be understood as generalizations of the original pp-wave solution studied in [22, 25, 97]. This work was extended by a more general study of multi-spin string solutions in [100], where the authors provided a detailed prescription for making direct comparisons with perturbative gauge theory. (For a more complete review of the development and current status of semiclassical string theory and the match-up with gauge theory, see [53] and references therein.) Early indications of integrability in the classical limit of the string theory emerged when it was shown that a certain configuration of the GS superstring action on $AdS_5 \times S^5$ admits an infinite set of classically conserved non-local charges, and may therefore be an integrable theory itself [59] (see also [40] for a reduction to the pp-wave system). The gauge theory analogue of this non-local symmetry was studied in [47, 48], where a direct connection with the string analysis was made to one-loop order in λ .

Various subtleties surrounding studies of the non-local (or Yangian) algebra arise at higher loops, and further work is certainly warranted.

In addition to the sector of non-local charges, however, integrable systems typically admit an infinite tower of local, mutually commuting charges, each of which is diagonalized by a set of Bethe equations [67, 101]. The presence of such a sector of hidden, classically conserved bosonic charges in the string theory was pointed out in [51]. Moreover, in accordance with the expectations of AdS/CFT duality, various studies have been successful in matching hidden local charges in the classical string theory to corresponding quantities in the quantum spin chain formulation of $\mathcal{N} = 4$ SYM theory. In [41], for example, Arutyunov and Staudacher constructed an infinite series of conserved local charges in the bosonic string theory by solving the Bäcklund equations associated with certain extended classical solutions of the $O(6)$ string sigma model. The local charges generated by the Bäcklund transformations were then matched to corresponding conserved charges obtained from an integrable quantum spin chain on the gauge theory side. In fact, they were able to demonstrate agreement between both sides of the duality for the entire infinite tower of local commuting charges. This study was extended in [42, 43], where it was shown that a general class of rotating classical string solutions can be mapped to solutions of a Neumann (or Neumann-Rosochatius) integrable system. More recently, a class of three-spin classical string solutions was shown in [102] to generate hidden local charges (again via Bäcklund transformations) that match their gauge theory counterparts to one-loop order. (For a thorough review of the match-up of semiclassical string integrable structures with corresponding structures in the gauge theory, see also [35, 53].)

The mapping between string and gauge theory integrable structures was studied from a somewhat different perspective in [50], where it was shown that the generator of local, classically conserved currents in the string theory is related in certain sectors to a particular Riemann-Hilbert problem that is reproduced precisely by the gauge theory integrable structure at one and two loops in λ . An analogous treat-

ment of the corresponding Riemann-Hilbert problem in non-compact sectors of the gauge/string duality was carried out in [49], and an extension of these studies to $\mathfrak{so}(6)$ and $\mathfrak{su}(2, 2)$ sectors was recently achieved in [46] and [103], respectively. The structure of the higher-loop Riemann-Hilbert problem descending from the classical string theory and its relationship with the corresponding gauge theory problem was used in conjunction with the long-range gauge theory Bethe ansatz of [34] to develop an ansatz that, albeit conjecturally, is purported to interpolate between the classical and quantum regimes of the string theory [44]. Although this proposal is not a proof of quantum integrability on the string side, it was demonstrated in [60] that the quantized string theory in the near-pp-wave limit yields a general multi-impurity spectrum that matches the string Bethe ansatz spectrum of [44]. The intricacy of this match-up is quite remarkable, and stands as strong evidence that this ansatz is correct for the string theory, at least to $O(1/J)$ in the large angular momentum (or background curvature) expansion. Furthermore, the proposed string Bethe equations can accommodate the strong-coupling $\lambda^{1/4}$ scaling behavior predicted in [104]. The spin chain theory implied by these Bethe equations, however, appears to disagree with that of the gauge theory, even at weak coupling [105].

Although the Bethe equations of [44] reproduce several predictions of the string theory in a highly nontrivial way, a direct test of quantum integrability (beyond tree level) in the string theory is still needed: this is the intent of the present chapter. Early steps in this direction were taken in [71], where the presence of a conserved local charge responsible for a certain parity degeneracy in the near-pp-wave string spectrum is examined at sixth-order in field fluctuations, or at $O(1/J^2)$ in the large- J expansion. Various subtleties of the analysis (possibly involving the proper renormalization of the theory at $O(1/J^2)$ in the expansion) make it difficult to reach any concrete conclusions, however. In this chapter we take a more immediate approach, relying primarily on a Lax representation of the classical string sigma model and studying a semiclassical expansion about certain point-like solitonic solutions. The goal is to establish the existence of a series of conserved, mutually commuting charges

in the string theory that can be quantized and studied using first-order perturbation theory. By aligning field fluctuations with the finite-radius curvature expansion studied in Chapters 3 and 4, we are able to study quantum corrections to quartic order, or to one loop beyond tree level. We show directly that several of the low-lying hidden charges in the series are conserved by the quantum theory to this order in the expansion, and we propose a method for matching specific eigenvalues of these charges to corresponding spectral quantities in the gauge theory.

The chapter is organized as follows: In Section 6.1 we briefly review the procedure for string quantization in the near-pp-wave limit developed in [26, 27], with the particular goal of demonstrating how background curvature corrections to the pp-wave theory can be interpreted as quantum corrections in a particular semiclassical expansion about point-like classical string solutions. In Section 6.2 we show how a Lax representation of the $O(4, 2) \times O(6)$ nonlinear sigma model can be modified to encode the string dynamics to the order of interest in this semiclassical expansion. We then generate a series of hidden local charges by expanding a perturbed monodromy matrix of the Lax representation in powers of the spectral parameter. In Section 6.3 we compute the eigenvalues of these charges in certain protected subsectors of the theory in the space of two-impurity string states. The resulting spectra are then compared on the S^5 subspace with those of corresponding charges descending from the $\mathfrak{su}(2)$ integrable sector of the gauge theory. We provide a prescription for matching the spectra of local charges on both sides of the duality, and carry out this matching procedure to eighth order in the spectral parameter. To the extent that they can be compared reliably, the gauge and string theory predictions are shown to match to this order (and presumably continue to agree at higher orders). We are thus led to believe that the integrable structure of the classical string theory survives quantization, at least to the first subleading order in field fluctuations beyond tree level.

6.1 Semiclassical string quantization in $AdS_5 \times S^5$

Most of the literature comparing semiclassical bosonic string theory in $AdS_5 \times S^5$ to corresponding sectors of gauge theory operators has focused on classical extended string solutions to the worldsheet sigma model in either “folded” or “circular” configurations, where certain components of the string angular momentum (i.e., certain charges of the Cartan subalgebra of the global symmetry group) are taken to be large (see, e.g., [42, 99, 100]). The latter amounts to choosing a so-called spinning ansatz for the string configuration [42, 43, 51, 53, 99, 100], and solutions endowed with such an ansatz can be identified with periodic solutions of the Neumann (or Neumann-Rosochatius) integrable system. The standard bosonic worldsheet action is usually chosen with flat worldsheet metric so that it is easily rewritten in terms of \mathbf{R}^6 embedding coordinates and identified with an $O(4, 2) \times O(6)$ sigma model. In the present study we will modify this treatment to allow for curvature corrections to the worldsheet metric, a complication that we are forced to confront when moving beyond tree level in lightcone gauge [26, 27].

We begin with a particular form of the $AdS_5 \times S^5$ target space metric, chosen originally in Chapters 3 and 4 for the fact that it admits a simple form for the spin connection:

$$ds^2_{AdS_5 \times S^5} = \widehat{R}^2 \left[- \left(\frac{1 + \frac{1}{4}z^2}{1 - \frac{1}{4}z^2} \right)^2 dt^2 + \left(\frac{1 - \frac{1}{4}y^2}{1 + \frac{1}{4}y^2} \right)^2 d\phi^2 + \frac{dz_k dz_k}{(1 - \frac{1}{4}z^2)^2} + \frac{dy_{k'} dy_{k'}}{(1 + \frac{1}{4}y^2)^2} \right]. \quad (6.1.1)$$

While we will not address fermions in this study, we will eventually return to the crucial issues of supersymmetry, and the metric choice in eqn. (6.1.1) will undoubtedly simplify further investigations. By defining

$$\cosh \rho \equiv \frac{1 + \frac{1}{4}z^2}{1 - \frac{1}{4}z^2}, \quad \cos \theta \equiv \frac{1 - \frac{1}{4}y^2}{1 + \frac{1}{4}y^2}, \quad (6.1.2)$$

we may write the $\mathbf{R}^6 \times \mathbf{R}^6$ embedding coordinates of AdS_5 and S^5 as

$$\begin{aligned} Z_k &= \sinh \rho \frac{z_k}{||z||} , & Z_0 + iZ_5 &= \cosh \rho e^{it} , \\ Y_{k'} &= \sin \theta \frac{y_{k'}}{||y||} , & Y_5 + iY_6 &= \cos \theta e^{i\phi} , \end{aligned} \quad (6.1.3)$$

with $||z|| \equiv \sqrt{z_k z_k}$. The coordinates Z_P , with $P, Q = 0, \dots, 5$, parameterize AdS_5 and are contracted over repeated indices using the metric $\eta_{PQ} = (-1, 1, 1, 1, 1, -1)$. The coordinates Y_M , with $M, N = 1, \dots, 6$, encode the S^5 geometry, and are contracted with a Euclidean metric.

Decomposing the theory into AdS_5 and S^5 subspaces, the usual conformal-gauge worldsheet action

$$S = - \int d^2 \sigma h^{ab} G_{\mu\nu} \partial_a x^\mu \partial_b x^\nu \quad (6.1.4)$$

can be written as

$$\begin{aligned} S &= \int d^2 \sigma (\mathcal{L}_{AdS_5} + \mathcal{L}_{S^5}) , \\ \mathcal{L}_{AdS_5} &= -\frac{1}{2} h^{ab} \eta_{PQ} \partial_a Z_P \partial_b Z_Q + \frac{\tilde{\varphi}}{2} (\eta_{PQ} Z_P Z_Q + 1) , \end{aligned} \quad (6.1.5)$$

$$\mathcal{L}_{S^5} = -\frac{1}{2} h^{ab} \partial_a Y_M \partial_b Y_M + \frac{\varphi}{2} (Y_M Y_M - 1) . \quad (6.1.6)$$

The quantities φ and $\tilde{\varphi}$ act as Lagrange multipliers in the action, enforcing the following conditions:¹

$$\eta_{PQ} Z_P Z_Q = -1 , \quad Y_M Y_M = 1 . \quad (6.1.7)$$

The action in eqn. (6.1.4) must also be supplemented by the standard conformal gauge constraints, and the worldsheet metric h^{ab} (the worldsheet indices run over $a, b \in \tau, \sigma$) will be allowed to acquire curvature corrections in accordance with these

¹Note that, in general, φ and $\tilde{\varphi}$ will depend on dynamical variables. We thank Arkady Tseytlin for clarification on this point.

constraints.

We wish to study a semiclassical expansion about the following classical point-like (or “BMN-like”) solutions to the sigma model equations of motion:

$$t = \phi = p_- \tau, \quad z_k = y_k = 0. \quad (6.1.8)$$

The expansion is defined in terms of quantum field fluctuations according to the following rescaling prescription:

$$t \rightarrow x^+, \quad \phi \rightarrow x^+ + \frac{x^-}{\sqrt{\xi}}, \quad z_k \rightarrow \frac{z_k}{\xi^{1/4}}, \quad y_k \rightarrow \frac{y_k}{\xi^{1/4}}. \quad (6.1.9)$$

(A similar but notably different choice was made in [99].) This particular choice of lightcone coordinates will allow us to maintain a constant momentum distribution on the worldsheet. Additionally, as noted in Chapter 3, it will have the effect of eliminating all normal-ordering ambiguities from the resulting worldsheet theory, an outcome that is particularly desirable in the present study. Furthermore, we note that if we identify $\xi \equiv \hat{R}^4$, the proposed expansion about the classical solution in eqn. (6.1.8) is identical to the large-radius curvature expansion about the pp-wave limit of $AdS_5 \times S^5$ studied above [26, 27, 38]. In other words, we have chosen a perturbation to the classical point-like string geodesic that reproduces the target-space curvature perturbation to the pp-wave limit. The background metric in eqn. (6.1.1) thus yields the following large- \hat{R} expansion:

$$\begin{aligned} ds^2 = & 2dx^+dx^- - (x^A)^2(dx^+)^2 + (dx^A)^2 \\ & + \frac{1}{\hat{R}^2} \left[-2y^2dx^+dx^- + \frac{1}{2}(y^4 - z^4)(dx^+)^2 + (dx^-)^2 + \frac{1}{2}z^2dz^2 - \frac{1}{2}y^2dy^2 \right] \\ & + O(\hat{R}^{-4}), \end{aligned} \quad (6.1.10)$$

where the pp-wave geometry emerges at leading order.

The details of quantizing the string Hamiltonian in this setting are given in Chap-

ter 3 (see also [24, 82–85] for further details), though we will briefly review the salient points here. The lightcone Hamiltonian H_{LC} is the generator of worldsheet time translations, and is defined in terms of the Lagrangian by

$$-H_{\text{LC}} = -p_+ = \delta\mathcal{L}/\delta\dot{x}^+ , \quad (6.1.11)$$

(or $\Delta - J$ in the language of BMN), and this variation is performed prior to any gauge fixing. The non-physical lightcone variables x^\pm are removed from the Hamiltonian by fixing lightcone gauge $x^+ = p_- \tau$ and replacing x^- with dynamical variables by enforcing the conformal gauge constraints

$$T_{ab} = \frac{\delta\mathcal{L}}{\delta h^{ab}} = 0 . \quad (6.1.12)$$

This procedure can be defined order-by-order in the large- \widehat{R} expansion. At leading order, for example, we obtain the following from eqn. (6.1.12):

$$\begin{aligned} \dot{x}^- &= \frac{p_-}{2}(x^A)^2 - \frac{1}{2p_-} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] + O(1/\widehat{R}^2) , \\ x'^- &= -\frac{1}{p_-} \dot{x}^A x'^A + O(1/\widehat{R}^2) . \end{aligned} \quad (6.1.13)$$

The conformal gauge constraints themselves are only consistent with the equations of motion if the worldsheet metric acquires curvature corrections (i.e., h departs from the flat metric $h = \text{diag}(-1, 1)$), which we express symbolically as \tilde{h}^{ab} according to

$$h = \begin{pmatrix} -1 + \tilde{h}^{\tau\tau}/\widehat{R}^2 & \tilde{h}^{\tau\sigma}/\widehat{R}^2 \\ \tilde{h}^{\tau\sigma}/\widehat{R}^2 & 1 + \tilde{h}^{\sigma\sigma}/\widehat{R}^2 \end{pmatrix} . \quad (6.1.14)$$

The requirement that $\det h = -1$ implies $\tilde{h}^{\tau\tau} = \tilde{h}^{\sigma\sigma}$ and, for future reference, the correction terms \tilde{h}^{ab} are given explicitly to the order of interest by

$$\tilde{h}^{\tau\tau} = \frac{1}{2}(z^2 - y^2) - \frac{1}{2p_-^2} \left[(\dot{x}^A)^2 + (x'^A)^2 \right] ,$$

$$\tilde{h}^{\tau\sigma} = \frac{1}{p_-^2} \dot{x}^A x'^A . \quad (6.1.15)$$

Finally, we note that the canonical momenta associated with the physical worldsheet excitations, defined by the variation $p_A = \delta\mathcal{L}/\delta x^A$, also acquire $O(1/\widehat{R}^2)$ corrections: consistent quantization requires that these corrections be taken into account. Expressed in terms of canonical variables, the final bosonic Hamiltonian takes the form

$$\begin{aligned} H_{\text{LC}} = & \frac{p_-}{2\widehat{R}^2} (x^A)^2 + \frac{1}{2p_- \widehat{R}^2} \left((p_A)^2 + (x'^A)^2 \right) \\ & + \frac{1}{\widehat{R}^4} \left\{ \frac{1}{4p_-} \left[z^2 (p_y^2 + y'^2 + 2z'^2) - y^2 (p_z^2 + z'^2 + 2y'^2) \right] + \frac{p_-}{8} [(x^A)^2]^2 \right. \\ & \left. - \frac{1}{8p_-^3} \left\{ [(p_A)^2]^2 + 2(p_A)^2 (x'^A)^2 + [(x^A)^2]^2 \right\} + \frac{1}{2p_-^3} (x'^A p_A)^2 \right\} + O(1/\widehat{R}^6) , \end{aligned} \quad (6.1.16)$$

where the pp-wave Hamiltonian emerges as expected at leading order. The lightcone momentum p_- is identified (via the AdS/CFT dictionary) with the modified 't Hooft parameter λ' according to

$$p_- = 1/\sqrt{\lambda'} = J/\sqrt{\lambda} . \quad (6.1.17)$$

From the point of view of the semiclassical analysis, we are working to two-loop order in quantum corrections. Since the quadratic theory can be quantized exactly, however, we can study the quartic interaction Hamiltonian using standard first-order perturbation theory. A detailed analysis of the resulting spectrum of this perturbation can be found above in Chapters 3, 4 and 5. In the course of those studies it was noticed that, analogous to the gauge theory closed sectors studied in [30, 31, 33, 69], certain sectors emerged from the string analysis that decouple from the remainder of the theory to all orders in λ' . One sector, which maps to the $\mathfrak{sl}(2)$ sector of the gauge theory, is diagonalized by bosonic string states excited in the AdS_5 subspace and forming symmetric-traceless irreps in spacetime indices. The corresponding sector of

symmetric-traceless S^5 string bosons maps to the closed $\mathfrak{su}(2)$ sector in the gauge theory. The block-diagonalization of these sectors in the string Hamiltonian will be an important tool in the present analysis: just as all higher hidden local charges in the gauge theory are simultaneously diagonalized by a single Bethe ansatz, all of the higher hidden charges descending from the string theory should be block-diagonalized by these particular string states as well.

6.2 Lax representation

The goal is to determine whether a ladder of higher local charges can be computed and quantized (albeit perturbatively), analogous to the existing treatment of the near-pp-wave Hamiltonian given in eqn. (6.1.16) above. To quartic order in the semiclassical expansion defined by eqn. (6.1.9), the difference between the string sigma model in eqns. (6.1.5, 6.1.6) and that of the $O(4, 2) \times O(6)$ sigma model, defined by

$$\begin{aligned}\mathcal{L}_{O(4,2)} &= -\frac{1}{2}\eta_{PQ}\partial_a Z_P\partial^a Z_Q + \frac{\tilde{\varphi}}{2}(\eta_{PQ}Z_P Z_Q + 1) , \\ \mathcal{L}_{O(6)} &= -\frac{1}{2}\partial_a Y_M\partial^a Y_M + \frac{\varphi}{2}(Y_M Y_M - 1) ,\end{aligned}\tag{6.2.1}$$

will essentially amount to an interaction perturbation due to curvature corrections to the worldsheet metric. We therefore find it useful to rely on a known Lax representation of the $O(4, 2) \times O(6)$ sigma model; this representation will define an unperturbed theory, and we will add perturbations by hand to recover the full interaction Hamiltonian in eqn. (6.1.16). (For a general introduction to the Lax methodology in integrable systems, the reader is referred to [101].) Since worldsheet curvature corrections only appear at $O(1/\hat{R}^2)$, the reduction to the $O(4, 2) \times O(6)$ sigma model at leading order in the expansion will be automatic.

For simplicity, we start from the four-dimensional Lax representation given for the $O(6)$ sigma model in [43] (see also [106] for details), and work only to leading order

in the semiclassical expansion. The complexified coordinates

$$\mathcal{Y}_1 = Y_1 + i Y_2 , \quad \mathcal{Y}_2 = Y_3 + i Y_4 , \quad \mathcal{Y}_3 = Y_5 + i Y_6 , \quad (6.2.2)$$

are used to form a unitary matrix S_{S^5}

$$S_{S^5} = \begin{pmatrix} 0 & \mathcal{Y}_1 & -\mathcal{Y}_2 & \bar{\mathcal{Y}}_3 \\ -\mathcal{Y}_1 & 0 & \mathcal{Y}_3 & \bar{\mathcal{Y}}_2 \\ \mathcal{Y}_2 & -\mathcal{Y}_3 & 0 & \bar{\mathcal{Y}}_1 \\ -\bar{\mathcal{Y}}_3 & -\bar{\mathcal{Y}}_2 & -\bar{\mathcal{Y}}_1 & 0 \end{pmatrix} , \quad (6.2.3)$$

in terms of which one may form the following $SU(4)$ -valued currents:

$$A_a = S_{S^5} \partial_a S_{S^5}^\dagger . \quad (6.2.4)$$

The equations of motion of the $O(6)$ sigma model

$$\partial_a \partial^a Y_M + \varphi Y_M = 0 \quad (6.2.5)$$

are then encoded by the auxiliary system of linear equations

$$(\partial_\sigma - U)X = (\partial_\tau - V)X = 0 , \quad (6.2.6)$$

where the Lax pair U and V are defined by

$$U = \frac{1}{1+\gamma} A_- - \frac{1}{1-\gamma} A_+ , \quad V = -\frac{1}{1+\gamma} A_- - \frac{1}{1-\gamma} A_+ . \quad (6.2.7)$$

The constant γ is a free spectral parameter, and A_\pm are defined by $A_\pm \equiv \frac{1}{2}(A_\tau \pm A_\sigma)$. Note that on the $SO(4)$ subspace spanned by $y_{k'}$, eqn. (6.2.5) reduces to the pp-wave equations of motion on S^5 :

$$\ddot{y}_{k'} - y''_{k'} + p_-^2 y_{k'} = 0 . \quad (6.2.8)$$

The utility of the Lax representation arises from the fact that U and V may be considered as local connection coefficients, and a consistency equation for the auxiliary linear problem can be reinterpreted as a flatness condition for the (U, V) -connection:

$$\partial_\tau U - \partial_\sigma V + [U, V] = 0 . \quad (6.2.9)$$

Parallel transport along this flat connection is defined by the path-ordered exponent

$$\Omega_C(\gamma) = \mathcal{P} \exp \int_C (U d\sigma + V d\tau) , \quad (6.2.10)$$

where \mathcal{C} is some contour in \mathbf{R}^2 . Restricting to transport along the contour defined by $\tau = \tau_0$ and $0 \leq \sigma \leq 2\pi$ yields a monodromy matrix:

$$T(2\pi, \gamma) = \mathcal{P} \exp \int_0^{2\pi} d\sigma U . \quad (6.2.11)$$

The flatness condition in eqn. (6.2.9) admits an infinite number of conservation laws, which translates to the fact that the trace of the monodromy matrix yields an infinite tower of local, mutually commuting charges $\widehat{Q}_n^{S^5}$ when expanded in powers of the spectral index about the poles of U ($\gamma = \pm 1$, in this case):²

$$\mathrm{tr} T(2\pi, \gamma) = \sum_n \gamma^n \widehat{Q}_n^{S^5} . \quad (6.2.12)$$

The first nonvanishing charge $\widehat{Q}_2^{S^5}$, for example, is the Hamiltonian of the theory (on the S^5 subspace).

Moving beyond leading order in the semiclassical expansion, the essential difference between the $O(6)$ sigma model defined in eqn. (6.2.1) and the string action given in eqn. (6.1.6) is, as noted above, that worldsheet indices are contracted in the latter case with a non-flat worldsheet metric. Keeping the components of h^{ab} explicit, the

²In general, an expansion around some γ that is finitely displaced from a singularity of U will yield combinations of local and non-local quantities. One is of course free to redefine γ such that the expansion about $\gamma = 0$ in eqn. (6.2.12) is local.

lightcone Hamiltonian derived from the string sigma model in eqn. (6.1.6) appears at leading order as

$$H_{\text{LC}}^{S^5} = -\frac{1}{2p_- \widehat{R}^2} \left[h^{\tau\tau} (p_-^2 y^2 + y'^2 + \dot{y}^2) + 2h^{\tau\sigma} \dot{y} \cdot y' \right] + O(1/\widehat{R}^4) , \quad (6.2.13)$$

where $h^{\tau\tau} = -1 + \tilde{h}^{\tau\tau}/\widehat{R}^2$ and $h^{\tau\sigma} = \tilde{h}^{\tau\sigma}/\widehat{R}^2$. The prescription will be to find a perturbation to the (U, V) -connection such that the Hamiltonian in eqn. (6.2.13) emerges in an appropriate limit from the charge $\widehat{Q}_2^{S^5}$ defined by eqn. (6.2.12). Such a perturbation is achieved by transforming the U matrix according to

$$U \rightarrow U = \frac{1}{1+\gamma} (1 + u_-/\widehat{R}^2) A_- - \frac{1}{1-\gamma} (1 + u_+/\widehat{R}^2) A_+ , \quad (6.2.14)$$

where u_{\pm} are given by

$$u_{\pm} \equiv -\frac{1}{2} \tilde{h}^{\tau\tau} \mp \frac{1}{3} \tilde{h}^{\tau\sigma} . \quad (6.2.15)$$

These perturbations should be treated as constants, to be replaced in the end with dynamical variables by fixing conformal gauge according to eqn. (6.1.12). The remaining quartic perturbations to the pp-wave theory will be naturally encoded in the semiclassical expansion of the underlying $O(6)$ (likewise, $O(4, 2)$) sigma model. The matrix V can be transformed in a similar way:

$$V \rightarrow V = -\frac{1}{1+\gamma} (1 + v_-/\widehat{R}^2) A_- - \frac{1}{1-\gamma} (1 + v_+/\widehat{R}^2) A_+ , \quad (6.2.16)$$

where v_{\pm} may be chosen such that the perturbed Lax pair satisfies the flatness condition in eqn. (6.2.9). Given that the intent is simply to determine whether the higher local charges generated by the perturbed monodromy matrix are conserved when quantum fluctuations are included, fixing V to satisfy the flatness condition is not really necessary: the complicated formulas for v_{\pm} that do satisfy eqn. (6.2.9) will therefore not be needed.

The perturbation in eqn. (6.2.14) can be obtained by a slightly different method. When the path-ordered exponent defining the monodromy matrix is expanded, it can be seen that all odd products of the Lax matrix U will not contribute to the final expression. By replacing all even products of U according to the rule

$$U(\sigma_1)U(\sigma_2) \rightarrow \frac{1}{(\gamma^2 - 1)^2} \left[h^{\sigma\sigma} A_\sigma(\sigma_1)A_\sigma(\sigma_2) - \gamma^2 h^{\sigma\tau} A_\sigma(\sigma_1)A_\tau(\sigma_2) - \gamma^2 h^{\tau\sigma} A_\tau(\sigma_1)A_\sigma(\sigma_2) - \gamma^2 h^{\tau\tau} A_\tau(\sigma_1)A_\tau(\sigma_2) \right], \quad (6.2.17)$$

the Hamiltonian in eqn. (6.2.13) is again obtained at leading order in the expansion. Computationally, this latter method seems to be much more efficient, and we will use eqn. (6.2.17) in what follows. At leading order in the $1/\widehat{R}$ expansion, the first nonvanishing integral of motion descending from the monodromy matrix is thereby found to be

$$Q_2^{S^5} = \frac{4\pi}{\widehat{R}^2} \int_0^{2\pi} d\sigma \left[h^{\tau\tau} (p_-^2 y^2 + y'^2 + \dot{y}^2) + 2h^{\tau\sigma} \dot{y} \cdot y' \right] + O(1/\widehat{R}^4), \quad (6.2.18)$$

which, by construction, matches the desired structure in eqn. (6.2.13).

The same construction may be carried out for the AdS_5 system. In fact, to make matters simple, we may borrow the Lax structure of the $O(6)$ model defined in eqns. (6.2.3-6.2.7), replacing the $O(6)$ coordinates in eqn. (6.2.2) with the following Euclideanized $O(4, 2)$ complex embedding coordinates:

$$\mathcal{Z}_1 = Z_1 + i Z_2, \quad \mathcal{Z}_2 = Z_3 + i Z_4, \quad \mathcal{Z}_3 = i Z_0 - Z_5. \quad (6.2.19)$$

In this case, however, the Lax matrix S_{AdS_5} will obey $S_{AdS_5}^\dagger S_{AdS_5} = -1$. Otherwise, the analysis above applies to the AdS_5 sector by direct analogy: expanding the perturbed $O(4, 2)$ monodromy matrix in the spectral parameter yields a set of charges labeled by $\widehat{Q}_n^{AdS_5}$. The local charges for the entire theory are then given by

$$\widehat{Q}_n \equiv \widehat{Q}_n^{S^5} - \widehat{Q}_n^{AdS_5}. \quad (6.2.20)$$

The corresponding currents will be labeled by \mathcal{Q}_n .

It turns out that the expansion in the spectral parameter γ is arranged such that the path-ordered exponent defining the monodromy matrix can be computed explicitly to a given order in γ by evaluating only a finite number of worldsheet integrals. The procedure for extracting local, canonically quantized currents is then completely analogous to that followed in computing the lightcone Hamiltonian described above. All gauge fixing is done after the currents are evaluated, all occurrences of x^- are replaced with dynamical variables by solving the conformal gauge constraints, and worldsheet metric corrections \tilde{h}^{ab} are evaluated according to eqns. (6.1.15) above. We note, however, that previous studies involving the matching of integrable structures between gauge and string theory have found it necessary to invoke certain redefinitions of γ to obtain agreement [34, 50]. It would be straightforward to allow for rather general redefinitions of the spectral parameter in the present calculation. When we turn to computing spectra and comparing with gauge theory, however, such redefinitions can lead to unwanted ambiguity. We will therefore be primarily interested in finding ratios of eigenvalue coefficients for which arbitrary redefinitions of γ are irrelevant, and for simplicity we will simply retain the original definition of γ given by eqn. (6.2.7) above.

As previously noted, the first current \mathcal{Q}_1 defined by eqn. (6.2.12) vanishes. In fact, all \mathcal{Q}_n vanish for odd values of n , and this property of the integrable structure is mirrored on the gauge theory side. The first nonvanishing current emerging from the monodromy matrix is given by

$$\begin{aligned} \mathcal{Q}_2 = & \frac{4\pi}{\widehat{R}^2} \left((\dot{x}^A)^2 + (x'^A)^2 + p_-^2 (x^A)^2 \right) \\ & + \frac{\pi}{\widehat{R}^4} \left\{ 2z^2 \left[y'^2 + 2z'^2 - \dot{y}^2 \right] - 2y^2 \left[z'^2 + 2y'^2 - \dot{z}^2 \right] - \frac{4}{p_-^2} (\dot{x}^A x'^A)^2 \right. \\ & \left. + \frac{1}{p_-^2} \left[3(\dot{x}^A)^2 - (x'^A)^2 \right] \left[(\dot{x}^A)^2 + (x'^A)^2 \right] + p_-^2 [(x^A)^2]^2 \right\} + O(1/\widehat{R}^6) . \end{aligned} \tag{6.2.21}$$

The leading-order term is the quadratic pp-wave Hamiltonian, as expected, and the perturbation is strictly quartic in field fluctuations. All occurrences of x^- and all curvature corrections to the worldsheet metric \tilde{h}^{ab} have been replaced with physical variables as described above. The final step is to express eqn. (6.2.21) in terms of canonically conjugate variables determined by directly varying the Lagrangian in eqn. (6.1.4). We obtain

$$\begin{aligned} \mathcal{Q}_2 = & \frac{4\pi}{\widehat{R}^2} \left(p_-^2 (x^A)^2 + (p_A)^2 + (x'^A)^2 \right) \\ & + \frac{\pi}{\widehat{R}^4} \left\{ 2 \left[-y^2 \left(p_z^2 + z'^2 + 2y'^2 \right) + z^2 \left(p_y^2 + y'^2 + 2z'^2 \right) \right] + p_-^2 [(x^A)^2]^2 \right. \\ & \left. - \frac{1}{p_-^2} \left\{ [(p_A)^2]^2 + 2(p_A)^2 (x'^A)^2 + [(x'^A)^2]^2 \right\} + \frac{4}{p_-^2} (x'^A p_A)^2 \right\} + O(1/\widehat{R}^6) . \end{aligned} \quad (6.2.22)$$

Comparing this with eqn. (6.1.16) above, we see that, to the order of interest,

$$\mathcal{Q}_2 = 8\pi p_- H_{LC} . \quad (6.2.23)$$

As expected, the perturbed monodromy matrix precisely reproduces the structure of the lightcone Hamiltonian to quartic order in the semiclassical expansion. (Note that \mathcal{Q}_2 is only expected to be identified with the lightcone Hamiltonian up to an overall constant.)

Computationally, the expansion of the monodromy matrix becomes increasingly time consuming at higher orders in the spectral index. The situation can be mitigated to some extent by projecting the theory onto AdS_5 or S^5 excitations, eliminating all interaction terms (from the quartic perturbation) that mix fluctuations from both subspaces. We will eventually want to compute eigenvalue spectra in the block-diagonal subsectors discussed above (which require such a projection), so this maneuver will not affect the outcome.

The next nonvanishing S^5 current in the series is given by

$$\begin{aligned}
\mathcal{Q}_4^{S^5} = & \frac{8\pi}{3\widehat{R}^2} (3 - \pi^2 p_-^2) (p_-^2 y^2 + p_y^2 + y'^2) \\
& + \frac{2\pi}{3p_-^2 \widehat{R}^4} \left\{ -3(p_y^2 - 2p_y \cdot y' + y'^2)(p_y^2 + 2p_y \cdot y' + y'^2) \right. \\
& - \pi^2 p_-^2 [4(p_y \cdot y')^2 + (p_y^2 + y'^2)^2] - 12p_-^2 y'^2 y^2 \\
& \left. - p_-^4 y^2 [4\pi^2 p_y^2 - 3y^2] - 3\pi^2 p_-^6 (y^2)^2 \right\} + O(1/\widehat{R}^6) . \quad (6.2.24)
\end{aligned}$$

Although the quadratic interaction of $\mathcal{Q}_4^{S^5}$ is proportional to the pp-wave Hamiltonian on the S^5 , the structure of the perturbing quartic interaction differs from that obtained for \mathcal{Q}_2 . The corresponding AdS_5 current takes the form

$$\begin{aligned}
\mathcal{Q}_4^{AdS_5} = & \frac{8\pi}{3\widehat{R}^2} (3 - \pi^2 p_-^2) (p_-^2 z^2 + p_z^2 + z'^2) \\
& + \frac{2\pi}{3p_-^2 \widehat{R}^4} \left\{ -3(p_z^2 - 2p_z \cdot z' + z'^2)(p_z^2 + 2p_z \cdot z' + z'^2) \right. \\
& - \pi^2 p_-^2 [4(p_z \cdot z')^2 + (p_z^2 + z'^2)^2] + 12p_-^2 z'^2 z^2 \\
& \left. + p_-^4 z^2 [-4\pi^2 z'^2 + 3z^2] + \pi^2 p_-^6 (z^2)^2 \right\} + O(1/\widehat{R}^6) , \quad (6.2.25)
\end{aligned}$$

where the quadratic sector is again proportional to the pp-wave Hamiltonian, projected in this case onto the AdS_5 subspace. Continuing on to sixth order in the spectral index, we find the S^5 current

$$\begin{aligned}
\mathcal{Q}_6^{S^5} = & \frac{1}{15\widehat{R}^2} \left\{ 4\pi [45 - 40\pi^2 p_-^2 + 2\pi^4 p_-^4] (p_-^2 y^2 + p_y^2 + y'^2) \right\} \\
& + \frac{\pi}{15p_-^2 \widehat{R}^4} \left\{ -45(p_y^2 - 2p_y \cdot y' + y'^2)(p_y^2 + 2p_y \cdot y' + y'^2) \right. \\
& - 20p_-^2 [2\pi^2 (4(p_y \cdot y')^2 + (p_y^2 + y'^2)^2) + 9y'^2 y^2] \\
& + p_-^4 [2\pi^4 (4(p_y \cdot y')^2 + 3(p_y^2 + y'^2)^2) - 160\pi^2 p_y^2 y^2 + 45(y^2)^2] \\
& \left. + 8\pi^2 p_-^6 y^2 [(2\pi^2 p_y^2 + \pi^2 y'^2) - 15y^2] + 10\pi^4 p_-^8 (y^2)^2 \right\} + O(1/\widehat{R}^6) . \quad (6.2.26)
\end{aligned}$$

The quadratic piece of $\mathcal{Q}_6^{S^5}$ is again identical in structure to the pp-wave Hamiltonian.

The analogous current in the AdS_5 subspace is arranged in a similar fashion:

$$\begin{aligned}
\mathcal{Q}_6^{AdS_5} = & \frac{1}{15\widehat{R}^2} \left\{ 4\pi \left[45 - 40\pi^2 p_-^2 + 2\pi^4 p_-^4 \right] \left(p_-^2 z^2 + p_z^2 + z'^2 \right) \right\} \\
& + \frac{\pi}{15p_-^2 \widehat{R}^4} \left\{ -45(p_z^2 - 2p_z \cdot z' + z'^2)(p_z^2 + 2p_z \cdot z' + z'^2) \right. \\
& - 20p_-^2 \left[2\pi^2 \left(4(p_z \cdot z')^2 + (p_z^2 + z'^2)^2 \right) - 9z'^2 z^2 \right] \\
& + p_-^4 \left[2\pi^4 \left(4(p_z \cdot z')^2 + 3(p_z^2 + z'^2)^2 \right) - 160\pi^2 z'^2 z^2 + 45(z^2)^2 \right] \\
& \left. + 8\pi^2 p_-^6 y^2 (\pi^2 z'^2 + 5z^2) - 6\pi^4 p_-^8 (z^2)^2 \right\} + O(1/\widehat{R}^6) . \tag{6.2.27}
\end{aligned}$$

While we will not present explicit formulas for the resulting currents, it is easy to carry this out to eighth order in γ .

Taken separately, each current can be viewed as a free pp-wave Hamiltonian plus a quartic interaction. This is particularly useful, as it allows us to quantize each charge exactly at leading order and express the perturbation in terms of free pp-wave oscillators. More explicitly, we quantize the quadratic sectors of these currents by expanding the fluctuation fields in their usual Fourier components:

$$\begin{aligned}
x^A(\sigma, \tau) &= \sum_{n=-\infty}^{\infty} x_n^A(\tau) e^{-ik_n \sigma} , \\
x_n^A(\tau) &= \frac{i}{\sqrt{2\omega_n}} \left(a_n^A e^{-i\omega_n \tau} + a_n^{A\dagger} e^{i\omega_n \tau} \right) . \tag{6.2.28}
\end{aligned}$$

The quadratic (pp-wave) equations of motion

$$\ddot{x}^A - x''^A + p_-^2 x^A = 0 \tag{6.2.29}$$

are satisfied by setting $k_n = n$ (integer), and $\omega_n = \sqrt{p_-^2 + k_n^2}$, where the operators a_n^A and $a_n^{A\dagger}$ obey the commutation relation $[a_m^A, a_n^{B\dagger}] = \delta_{mn} \delta^{AB}$.

In accordance with integrability, we expect that the local charges in eqns. (6.2.22–6.2.27) should all be mutually commuting. Expressed in terms of quantum raising

and lowering operators, we can check the commutators of the hidden local charges directly. To avoid mixing issues, we will need to select out closed subsectors of each charge that completely decouple from the remaining terms in the theory. We have already noted that the Hamiltonian \widehat{Q}_2 is known to be closed under AdS_5 and S^5 string states forming symmetric-traceless irreps in their spacetime indices. The equivalent gauge theory statement is that the dilatation generator is closed in certain $\mathfrak{sl}(2)$ and $\mathfrak{su}(2)$ projections. Since the complete tower of corresponding charges in the gauge theory (including the dilatation generator) can be diagonalized by a single set of $\mathfrak{sl}(2)$ or $\mathfrak{su}(2)$ Bethe equations, it is a reasonable guess that the full tower of local string charges decouples under corresponding projections. (A similar conjecture is made, for example, in [41, 50].) Following the treatment in Chapter 5, we therefore define the following AdS_5 oscillators

$$a_n = \frac{1}{\sqrt{2}} (a_n^j + ia_n^k) , \quad \bar{a}_n = \frac{1}{\sqrt{2}} (a_n^j - ia_n^k) , \quad (j \neq k) , \quad (6.2.30)$$

which satisfy the standard relations

$$[a_n, a_m^\dagger] = [\bar{a}_n, \bar{a}_m^\dagger] = \delta_{nm} , \quad [a_n, \bar{a}_m^\dagger] = [\bar{a}_n, a_m^\dagger] = 0 . \quad (6.2.31)$$

When restricted to these oscillators, the symmetric-traceless projection in the AdS_5 subspace is achieved by setting all $\bar{a}_n, \bar{a}_n^\dagger$ to zero (see [60] for details). A corresponding definition on the S^5 takes the form

$$a_n = \frac{1}{\sqrt{2}} (a_n^{j'} + ia_n^{k'}) , \quad \bar{a}_n = \frac{1}{\sqrt{2}} (a_n^{j'} - ia_n^{k'}) , \quad (j' \neq k') , \quad (6.2.32)$$

where the symmetric-traceless projection is again invoked by setting $\bar{a}_n, \bar{a}_n^\dagger$ to zero. In other words, we can test the commutativity of the local charges in the AdS_5 and S^5 symmetric-traceless projections by rewriting their oscillator expansions according to eqns. (6.2.30, 6.2.32) and setting all $\bar{a}_n, \bar{a}_n^\dagger$ to zero.

Since the currents are expanded to $O(1/\widehat{R}^4)$, we only require that the commu-

tators vanish to $O(1/\widehat{R}^6)$. This simplifies the problem somewhat, since we only need to compute commutators involving at most six oscillators. On the subspace of symmetric-traceless AdS_5 string states, we obtain

$$\left[\widehat{Q}_n^{AdS_5}, \widehat{Q}_m^{AdS_5} \right] = O(1/\widehat{R}^6) , \quad n, m \in 2, \dots, 8 . \quad (6.2.33)$$

The corresponding projection on the S^5 yields

$$\left[\widehat{Q}_n^{S^5}, \widehat{Q}_m^{S^5} \right] = O(1/\widehat{R}^6) , \quad n, m \in 2, \dots, 8 . \quad (6.2.34)$$

We therefore find evidence for the existence of a tower of mutually commuting charges (within these particular closed sectors) that are conserved perturbatively by the quantized theory.

6.3 Spectral comparison with gauge theory

Given the freedom involved in redefinitions of the spectral parameter, it may seem that any spectral agreement between the string charges computed above and corresponding quantities in the gauge theory would be rather arbitrary. We therefore seek a comparison of integrable structures on both sides of the duality that avoids this ambiguity. It turns out that such a test is indeed possible in the symmetric-traceless sector of S^5 excitations, which will map in the gauge theory to the closed $\mathfrak{su}(2)$ sector. We will further restrict ourselves to computing spectra associated with the following two-impurity string states:

$$a_q^{j'\dagger} a_{-q}^{k'\dagger} |J\rangle .$$

The analysis for three- or higher-impurity states would require an accounting of interactions between AdS_5 and S^5 string excitations; as noted above, however, this dramatically complicates the computational analysis. (We intend to return to the

question of higher-impurity string integrability in a future study.) The ground state $|J\rangle$ is understood to carry J units of angular momentum on the S^5 , and the two-impurity $SO(4)$ subspace above comprises a 16×16 -dimensional sub-block of the Hamiltonian. In addition, the mode indices (labeled here by q) of physical string states must sum to zero to satisfy the usual level-matching condition (the Virasoro constraint is understood to be satisfied by the leading-order solution to the equations of motion; any higher-order information contained in the T_{01} component of eqn. (6.1.12) is irrelevant).

To simplify the analysis, and for comparison with previous chapters, we will also rescale each of the charges computed above by a factor of \widehat{R}^2 :

$$\widehat{Q}_n \rightarrow \widehat{R}^2 \widehat{Q}_n . \quad (6.3.1)$$

The two-impurity matrix elements of the charge $\widehat{Q}_2^{S^5}$ are then given by:

$$\begin{aligned} \langle J | a_q^{a'} a_{-q}^{b'} (\widehat{Q}_2^{S^5}) a_{-q}^{c'\dagger} a_q^{d'\dagger} | J \rangle &= 16\pi\omega_q \delta^{a'd'} \delta^{b'c'} \\ &- \frac{8\pi q^2}{J\sqrt{\lambda'}\omega_q^2} \left[(3 + 2q^2\lambda) \delta^{a'd'} \delta^{b'c'} - \delta^{a'c'} \delta^{b'd'} + \delta^{a'b'} \delta^{c'd'} \right] + O(1/J^2) , \end{aligned} \quad (6.3.2)$$

The radius \widehat{R} has been replaced with the angular momentum J , and p_- has been replaced with $1/\sqrt{\lambda'}$ via

$$J/p_- = \widehat{R}^2 = \sqrt{\lambda'} . \quad (6.3.3)$$

As expected, contributions to the pp-wave limit of eqn. (6.3.2) all lie on the diagonal. Up to an overall factor, one may further check that the correction terms at $O(1/J)$ agree with those computed in Chapter 3, projected onto the S^5 subspace. The next

higher charges in the series yield matrix elements given by

$$\begin{aligned}
\langle J | a_q^{a'} a_{-q}^{b'} (\widehat{Q}_4^{S^5}) a_{-q}^{c'\dagger} a_q^{d'\dagger} | J \rangle &= \frac{32\pi\omega_q}{3\lambda'} (3\lambda' - \pi^2) \delta^{a'd'} \delta^{b'c'} \\
&\quad - \frac{16\pi}{3\lambda'^{5/2} \omega_q^2 J} \left\{ \left[\pi^2 (2 + q^2 \lambda') + 3q^2 \lambda'^2 (3 + 2q^2 \lambda') \right] \delta^{a'd'} \delta^{b'c'} \right. \\
&\quad \left. + \left[\pi^2 (2 + q^2 \lambda') - 3q^2 \lambda'^2 \right] \delta^{a'c'} \delta^{b'd'} \right. \\
&\quad \left. + q^2 \lambda' (3\lambda' - \pi^2) \delta^{a'b'} \delta^{c'd'} \right\} + O(1/J^2) , \tag{6.3.4}
\end{aligned}$$

$$\begin{aligned}
\langle J | a_q^{a'} a_{-q}^{b'} (\widehat{Q}_6^{S^5}) a_{-q}^{c'\dagger} a_q^{d'\dagger} | J \rangle &= \frac{16\pi}{15\omega_q \lambda'^2} (2\pi^4 - 40\pi^2 \lambda' + 45\lambda'^2) \omega_q^2 \delta^{a'd'} \delta^{b'c'} \\
&\quad + \frac{8\pi}{15\lambda'^{7/2} \omega_q^2} \left\{ \left[2\pi^4 (4 + q^2 \lambda' (5 + q^2 \lambda')) - 40\pi^2 \lambda' (2 + q^2 \lambda') \right. \right. \\
&\quad \left. \left. - 45q^2 \lambda'^3 (3 + 2q^2 \lambda') \right] \delta^{a'd'} \delta^{b'c'} + \left[2\pi^4 (4 + q^2 \lambda') - 40\pi^2 \lambda' (2 + q^2 \lambda') \right. \right. \\
&\quad \left. \left. + 45q^2 \lambda'^3 \right] \delta^{a'c'} \delta^{b'd'} + q^2 \lambda' \left[\lambda' (40\pi^2 - 45\lambda') - 2\pi^4 \right] \delta^{a'b'} \delta^{c'd'} \right\} + O(1/J^2) . \tag{6.3.5}
\end{aligned}$$

We will again project onto symmetric-traceless irreps of $SO(4) \times SO(4)$, transforming as $(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$ in an $SU(2)^2 \times SU(2)^2$ notation. Although it is not necessarily guaranteed that the symmetric-traceless states will diagonalize the higher charges \widehat{Q}_4 and \widehat{Q}_6 at quartic order, this can be checked directly at one-loop order in λ' by computing the eigenvectors of the charges above (the higher-loop version of this check is much more difficult because the above charges are no longer completely block diagonal under the $SO(4)$ projection, a fact that can be seen in the structure of \mathcal{Q}_2 above). The $\widehat{Q}_2^{S^5}$ eigenvalue between symmetric-traceless $(\mathbf{1}, \mathbf{1}; \mathbf{3}, \mathbf{3})$ S^5 states (denoted by $Q_2^{S^5}$) is then found to be

$$Q_2^{S^5} = 16\pi \left(\omega_q - \frac{q^2 \sqrt{\lambda'}}{J} \right) + O(1/J^2) . \tag{6.3.6}$$

Up to an overall constant, this is just the two-impurity energy shift computed in Chapter 3. The corresponding eigenvalues of the higher charges \widehat{Q}_4 and \widehat{Q}_6 can be

computed in an analogous fashion:

$$\begin{aligned}
Q_4^{S^5} &= \frac{32\pi}{3} \left\{ \frac{\omega_q}{\lambda'} [3\lambda' - \pi^2] \right. \\
&\quad \left. - \frac{\pi}{\lambda'^{5/2} \omega_q^2 J} [\pi^2(2 + q^2 \lambda') + 3q^2 \lambda'^3 \omega_q^2] \right\} + O(1/J^2) , \\
Q_6^{S^5} &= \frac{16\pi}{15} \left\{ \frac{\omega_q}{\lambda'^2} (2\pi^4 - 40\pi^2 \lambda' + 45\lambda'^2) - \frac{1}{\omega_q^2 \lambda'^{7/2} J} [40\pi^2 \lambda' (2 + q^2 \lambda') \right. \\
&\quad \left. + 45q^2 \lambda'^4 \omega_q^2 - 2\pi^4 (4 + q^2 \lambda' (3 + q^2 \lambda'))] \right\} + O(1/J^2) . \tag{6.3.7}
\end{aligned}$$

Similar formulas can be extracted for the AdS_5 charges $Q_2^{AdS_5}$, $Q_4^{AdS_5}$ and $Q_6^{AdS_5}$, which are diagonalized by symmetric-traceless $(\mathbf{3}, \mathbf{3}; \mathbf{1}, \mathbf{1})$ string states excited in the AdS_5 subspace. Though we have not given explicit formulas, it is also straightforward to obtain the corresponding eigenvalues for $Q_8^{AdS_5}$ and $Q_8^{S^5}$.

By modifying the Inozemtsev spin chain of [37] to exhibit higher-loop BMN scaling, Beisert, Dippel and Staudacher were able to formulate a long-range Bethe ansatz for the gauge theory in the closed $\mathfrak{su}(2)$ sector [34] (we will simply state their results here, referring the reader to [34] for further details). In essence, the Bethe ansatz encodes the interactions of pseudoparticle excitations on a spin lattice and, in terms of pseudoparticle momenta p_k , the ansatz given in [34] diagonalizes the entire tower of local gauge theory $\mathfrak{su}(2)$ charges. The eigenvalues of these charges, which we label here as D_n , are given by

$$D_n = \sum_{k=1}^I q_n(p_k) , \quad q_n(p) = \frac{2 \sin(\frac{p}{2}(n-1))}{n-1} \left(\frac{\sqrt{1 + 8g^2 \sin^2(p/2)} - 1}{2g^2 \sin^2(p/2)} \right)^{n-1} , \tag{6.3.8}$$

where $g^2 \equiv \lambda/8\pi^2$, and the index k runs over the total number I of pseudoparticle excitations (or R -charge impurities) on the spin lattice. These eigenvalues can then be expanded perturbatively in inverse powers of the gauge theory R -charge (R) by

approximating the pseudoparticle momenta p_k by the expansion

$$p_k = \sum_j \frac{f_j(n_k)}{R^{j/2}} , \quad (6.3.9)$$

where f_j are functions of the integer mode numbers n_k , determined by solving the Bethe equations explicitly to a given order in $1/R$.

In general, we wish to identify the local string charges with linear combinations of corresponding charges in the gauge theory. From eqn. (6.3.8), however, it is easy to see that as one moves up the ladder of higher charges in the gauge theory, the eigenvalues D_n of these charges have leading contributions at higher and higher powers of g^2/R^2 in the large- R , small- λ double-scaling expansion. This is puzzling because the string eigenvalues computed above do not exhibit similar properties. The difference in scaling behavior therefore motivates the following prescription for identifying the eigenvalues of the higher local charges on both sides of the correspondence:

$$Q_n - N = C \left(\frac{n}{2} D_n \right)^{2/n} . \quad (6.3.10)$$

N here counts the number of string worldsheet impurities and C is an arbitrary constant. Fractional powers of the gauge theory charges D_n are well defined in terms of the double-scaling expansion, so that the right-hand side of eqn. (6.3.10) is in fact just a linear combination of conserved quantities in the gauge theory.

A potential subtlety arises when matching Q_n and D_n in this fashion for $n > 2$ beyond one-loop order in λ . The problem is that, under the identification in eqn. (6.3.10), information from string energy eigenvalues at $O(1/J^2)$ and higher is required to completely characterize the higher-loop (in λ) coefficients of the gauge theory charges D_n . The essential reason for this is that the string loop expansion is in powers of the modified 't Hooft coupling, which, in terms of the gauge theory R -charge R , is

$$\lambda' = \lambda/J^2 = \lambda/R^2 . \quad (6.3.11)$$

In other words, under eqn. (6.3.10), it is impossible to disentangle higher-order $1/J$ contributions to the string charges Q_n from higher-order λ corrections to D_n . The prescription given in eqn. (6.3.10) therefore holds only to one-loop order in λ , where knowing the $1/J$ corrections in the string theory is sufficient.

Furthermore, since the local charges in the string and gauge theories are only identified up to an overall multiplicative constant, directly comparing the spectra of each theory is not especially rigorous. A convenient quantity to work with, however, is the ratio of the $O(1/J)$ eigenvalue correction to the pp-wave coefficient: at first-loop order in λ this ratio eliminates all ambiguity associated with overall constants and γ redefinitions, and thus provides a meaningful comparison with gauge theory. (The analogous quantity computed for charges in the AdS_5 subspace is not free from such ambiguities.) We therefore arrange the one-loop, two-impurity eigenvalues of local S^5 string theory charges according to

$$Q_n^{S^5} = 2 + q^2 \lambda' \left(\Lambda_{n,0} + \frac{\Lambda_{n,1}}{J} \right) + O(\lambda'^2) + O(1/J^2) , \quad (6.3.12)$$

where the numbers $\Lambda_{n,0}$ and $\Lambda_{n,1}$ characterize eigenvalue coefficients in the pp-wave limit and at $O(1/J)$, respectively, and q is the mode number associated with the two-impurity string states defined above. On the gauge theory side we make a similar arrangement:

$$\left(\frac{n}{2} D_n \right)^{2/n} = \frac{q^2 \lambda}{R^2} \left(\bar{\Lambda}_{n,0} + \frac{\bar{\Lambda}_{n,1}}{R} \right) + O(\lambda^2) + O(1/R^4) , \quad (6.3.13)$$

where the integer q is a mode number associated with the momenta of pseudoparticle excitations on the spin lattice (which, in turn, correspond to roots of the $\mathfrak{su}(2)$ Bethe equations). The R -charge R is understood to be identified with the string angular momentum J via eqn. (6.3.11).

The quantities $\Lambda_{2,0}$ and $\Lambda_{2,1}$ for the string Hamiltonian Q_2 can be computed from the eigenvalue formula in eqn. (6.3.6) (or, alternatively, retrieved from the two-

n	$\Lambda_{n,1}/\Lambda_{n,0}$	$\bar{\Lambda}_{n,1}/\bar{\Lambda}_{n,0}$
2	-2	-2
3	0	0
4	-2	-2
5	0	0
6	-2	-2
7	0	0
8	-2	-2

Table 6.1: Ratios of $O(1/J)$ (or $O(1/R)$) corrections to pp-wave/BMN coefficients in string and gauge theory local charges

impurity string results reported in Chapter 3). We find the following ratio:

$$\Lambda_{2,1}/\Lambda_{2,0} = -2 . \quad (6.3.14)$$

As shown in Chapter 3, this agrees with the corresponding gauge theory prediction at one-loop order in λ :

$$\bar{\Lambda}_{2,1}/\bar{\Lambda}_{2,0} = -2 . \quad (6.3.15)$$

The ratio of $O(1/J)$ eigenvalue corrections to pp-wave coefficients is in fact -2 for all of the nonvanishing string charges. Under the matching prescription in eqn. (6.3.10), this agrees with the gauge theory perfectly. (The odd charges vanish altogether on both sides of the correspondence.) We summarize the results of this comparison for the first eight charges in the series in Table 6.1. It would be satisfying to test this agreement at higher loop-orders in λ . The corresponding computation at two-loop order, however, would require evaluating the local string theory charges at $O(1/\hat{R}^6)$ in the semiclassical expansion, where several subtleties of perturbation theory (and, for that matter, lightcone quantization) would need to be addressed. This emphasizes the need to understand the quantum string theory at higher orders in the expansion away from the pp-wave limit.

6.4 Discussion

In this chapter we have provided evidence that an infinite tower of local, mutually commuting bosonic charges of type IIB string theory on $AdS_5 \times S^5$, known to exist in the classical theory, can also be identified in the quantum theory. In addition, we have provided a prescription for matching certain eigenvalues of these charges in a protected subsector of the string theory to corresponding eigenvalues in the closed $\mathfrak{su}(2)$ sector of the gauge theory. The fact that the spectra of local string charges computed here can only be matched to corresponding quantities in the gauge theory via the matching prescription in eqn. (6.3.10), however, indicates that the monodromy matrix used to derive the local string charges is substantially different from that which would give rise to the proposed quantum string Bethe ansatz of [44] (or, since they are equivalent at one-loop order, the corresponding $\mathfrak{su}(2)$ Bethe ansatz in the gauge theory). In other words, we expect that there is a Lax representation for the string sigma model that gives rise to hidden local charges that can be compared directly with the gauge theory, without having to take fractional powers or linear combinations.

There are a number of additional tests of integrability in the quantum string theory that, in the context of the present calculation, should be relatively straightforward. By computing the quartic interactions among fluctuations in the AdS_5 and S^5 subspaces for each of the higher local charges studied here, it would be easy, for example, to find the resulting spectra of three- or higher-impurity string states. Apart from the difficulty of actually computing the mixing interactions, this would provide a simple check on the methodology employed here. A more difficult problem would be to address whether the integrable structure of the string theory respects supersymmetry. By formulating a supersymmetric Lax representation that generates the complete interaction Hamiltonian computed in Chapter 3, one might be able to show that each of the higher local charges are individually supersymmetric, and a comparison with gauge theory could be carried out in the closed $\mathfrak{su}(1|1)$ sector studied in Chapters 4 and 5 (the corresponding sector of the string theory would be comprised of

symmetrized fermionic excitations in the $(\mathbf{3}, \mathbf{1}; \mathbf{3}, \mathbf{1})$ or $(\mathbf{1}, \mathbf{3}; \mathbf{1}, \mathbf{3})$ of $SO(4) \times SO(4)$.

Ultimately, the hope is that the arsenal of techniques associated with integrable systems can be employed to find an exact solution to the string formulation of large- N_c Yang-Mills theory. Alternatively, a proof that both sides of the duality are diagonalized by identical Bethe equations should be obtainable. At present, the major obstacle preventing such a proof is the disagreement between gauge and string theory at three-loop order in the 't Hooft coupling. The fact that the integrable systems of both theories seem to agree in certain limited cases, however, stands as strong evidence that they are likely to be equivalent.

Chapter 7

Conclusions and outlook

Since the advent of the BMN/pp-wave limit, the AdS/CFT correspondence has been subject to a new class of rigorous and detailed tests. These studies have not only provided novel verifications of the validity of Maldacena's conjecture, but they have given a much more detailed understanding of how holographic dualities are realized. In this dissertation we have focused on the correspondence between type IIB superstring theory on $AdS_5 \times S^5$ and $\mathcal{N} = 4$ SYM theory in four dimensions, widely viewed as the simplest and most striking example of AdS/CFT duality. Given the large number of symmetries on either side of the correspondence, this is, in many respects, the easiest system to study. It has been a longstanding problem, however, that string quantization is not well understood in the presence of a curved, RR background. Together with the difficulty of computing non-BPS operator dimensions in the strong-coupling limit of the gauge theory, the obstacles preventing a direct test of the proposal have been formidable. Of course, this situation changed dramatically when BMN discovered a large R -charge limit of the gauge theory that matched the pp-wave limit of the string theory, a limit that was shown by Metsaev to render the string theory exactly soluble. Their insights were a tremendous success because the match-up between the string and gauge theory in this limit marked the first direct comparison of string energy spectra with a corresponding set of anomalous dimensions in the gauge theory.

At this level of the analysis, however, one is limited to dealing with spectra on

either side of the duality that are highly degenerate. We demonstrated that this degeneracy can be lifted by including worldsheet interactions associated with background curvature corrections to the Penrose limit. Including such corrections is a difficult task, as it reintroduces the puzzle of quantizing the string theory in a curved background with RR flux. The essential point is that this vexing problem can be circumvented by treating the curvature corrections in a purely perturbative setting. The corrections to the spectrum can be controlled by the expansion in inverse (squared) powers of the scale radius ($1/\hat{R}^{2n}$) or, equivalently, in inverse powers of the S^5 angular momentum ($1/J$). We have focused in this dissertation on the leading set of these corrections, appearing at $O(1/J)$ in the expansion.

Chapter 3 was dedicated to computing the interaction Hamiltonian and analyzing the resulting spectrum in the Fock space of two-impurity string states, formed by acting with two raising operators on the ground state $|J\rangle$. The resulting spectrum was composed of 256 distinct states that sort themselves into a nine-level supermultiplet whose multiplicity structure matches that which is expected from the structure of $\mathcal{N} = 4$ supersymmetry. Furthermore, the string theory provides energies that are exact in the gauge theory coupling $\lambda = g_{\text{YM}}^2 N_c$. When compared with higher-order λ corrections to anomalous dimensions computed directly from the gauge theory, we found a perfect match at both one- and two-loop order, but this remarkable agreement with the gauge theory breaks down at third order.

By extending this analysis to the 4,096-dimensional Fock space of three-impurity string states, we were able to show in Chapter 4 that the string and gauge theories again agree in this perturbative setting at one- and two-loop order. The agreement again breaks down at three loops, however. When compared with the conjectured Bethe equations that provide anomalous dimension formulas in the gauge theory, this pattern was discovered in Chapter 5 to exist for the generic N -impurity case, albeit restricted to certain protected subsectors of the theory.

The BMN/pp-wave limit of the AdS/CFT correspondence provides an immensely powerful testing ground for holographic gauge/string dualities. We have provided

direct spectral comparisons of the gauge and string theories, along with convincing evidence that, in this setting, one can expect to see an impressive match-up that holds to two-loop order in the 't Hooft coupling, but breaks down at three loops.¹ Whether one can perform the similar analyses in the string theory at even higher orders in the $1/\widehat{R}$ expansion is still an open question. (Early steps in this direction are taken in [71].) It is unclear whether lightcone methods will be helpful in this context, because, dimensionally speaking, the theory becomes nonrenormalizable at $O(1/\widehat{R}^4)$ in the expansion. In addition, it still remains to be seen whether the diverse set of integrable structures underlying the duality will be useful for actually solving certain sectors of the gauge theory or the string theory.

In the end we stand to learn a great deal about the non-perturbative aspects of Yang-Mills theories by relying on their string theory counterparts. As noted above, the next landmark achievement will likely be proving the complete equivalence of some sector of IIB string theory on $AdS_5 \times S^5$ and $\mathcal{N} = 4$ SYM theory, perhaps at the level of the Bethe equations. For this to happen, however, we need to rectify the higher-loop disagreement. Given the huge amount of symmetry on both sides of the duality, it may be possible to reach the strong-coupling regime of the gauge theory and verify the form of the conjectured wrapping interactions therein. If current trends continue, we will undoubtedly uncover fascinating new realms of physics in the process.

¹This pattern has been found elsewhere, particularly in the study of the duality between semi-classical extended string configurations and corresponding sectors of the gauge theory (see also [53] for a general review of this program, and [55, 56, 107–111] for more recent developments). Extended string configurations typically give rise to additional conserved charges that provide an intuitive generalization of the BMN/pp-wave picture.

Appendix A

Notation and conventions

For convenience we record in this appendix the most common symbols used in the text. In the following list we collect quantities defined on the CFT side of the correspondence:

D	Operator dimension
R	$U(1)_R$ component of the $SU(4)$ R -symmetry (the R -charge); is mapped to and used interchangeably with the string theory angular momentum J (see below)
K	Naive dimension; counts the total number of fields in an operator, or the number of sites on the corresponding spin chain; maps to $J + I$ on the string side, where I is the impurity number
Δ	$D - R$; maps to P_+ on the string side
g_{YM}	Yang-Mills coupling
λ	$g_{\text{YM}}^2 N_c$, 't Hooft coupling
N_c	Rank of the Yang-Mills gauge group
$\overline{T}_K^{(+)}$	Symmetric-traceless, rank-two $SO(4)$ tensor operator in the $(2, K - 4, 2)$ $SU(4)$ irrep
$\overline{T}_K^{(-)}$	Antisymmetric, rank-two $SO(4)$ tensor operator in the $(0, K - 3, 2) + (2, K - 3, 0)$ of $SU(4)$
$\overline{T}_K^{(0)}$	Trace part of the set of rank-two $SO(4)$ tensor operators,

	in the $(0, K - 2, 0)$ of $SU(4)$
$D - K$	Anomalous dimension
Δ_0	$K - R$, counts the number of R -charge impurities (can be half-integer valued when fermionic impurities are present)
L	The level of a supermultiplet, reached by acting on a primary level with some SUSY generator L times
ϕ^{\square}	SYM scalars, rank-two antisymmetric $SU(4)$ tensor in the six-dimensional $(0, 1, 0)$ ($(\mathbf{2}, \mathbf{1})$ of $SL(2, C)$)
χ^{\square}_a	SYM gluino, rank-one $SU(4)$ tensor in the four-dimensional fundamental $(1, 0, 0)$ ($(\mathbf{1}, \mathbf{2})$ of $SL(2, C)$)
$\bar{\chi}^{\square}_{\dot{a}}$	SYM gluino, rank-three antisymmetric $SU(4)$ tensor in the four-dimensional antifundamental $(0, 0, 1)$
∇_{μ}	Spacetime covariant derivative
Z	Scalar field with $R = 1$: $\phi^{\square}_{\mathbf{2}}$
A, B	Scalar impurity fields with $R = 0$: $\phi^{\square}_{\mathbf{3}}$, $\phi^{\square}_{\mathbf{4}}$, $\phi^{\square}_{\mathbf{3}}$, $\phi^{\square}_{\mathbf{4}}$ (or simply ϕ^A, ϕ^B)
\bar{Z}	Scalar field with $R = -1$: $\phi^{\square}_{\mathbf{4}}$
$P_{i,j}$	Permutation operator that exchanges spins on the i^{th} and j^{th} lattice sites of a spin chain
$\{n_1, n_2, \dots\}$	Shorthand for the following series of permutation operators: $\sum_{k=1}^L P_{k+n_1, k+n_1+1} P_{k+n_2, k+n_2+1} \dots$
I	Spin chain impurity number (Chapter 2)
n_i	Mode number of the i^{th} pseudoparticle excitation on the spin lattice (Chapter 2)
b_j^{\dagger}, b_j	Position space raising and lowering operators for magnon excitations on the spin lattice
$\tilde{b}_p^{\dagger}, \tilde{b}_p$	Momentum space raising and lowering operators for magnon excitations
$ L\rangle$	Ground state of the length- L spin chain; corresponds to the BPS operator $\text{tr}(Z^L)$

$\bar{\Lambda}$	$O(1/R)$ shift of the anomalous dimension
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On the string side we use the following symbols:

g_s	String coupling, equal to g_{YM}^2
\widehat{R}	Curvature radius of $AdS_5 \times S^5$
λ'	Modified 't Hooft coupling, $g_{\text{YM}}^2 N_c / J^2$
J	Angular momentum of string states along an equatorial geodesic in the S^5 subspace; maps to $L - I$, or simply R , in the CFT
ω	String excitation energy
$ J\rangle$	The string ground state, carrying J units of angular momentum on the S^5
P_+	String lightcone Hamiltonian $\omega - J$, maps to Δ on the CFT side
L	Supermultiplet level of a string energy spectrum
a_n^\dagger, a_n	Bosonic raising and lowering operators with mode number n
b_n^\dagger, b_n	Fermionic raising and lowering operators with mode number n
z^k	$SO(4)$ vector in AdS_5
$y^{k'}$	$SO(4)$ vector in S^5
x^A	Vector in a transverse $SO(8)$ subspace of $AdS_5 \times S^5$
$G_{\mu\nu}$	Metric tensor of $AdS_5 \times S^5$
h_{ab}	Worldsheet metric tensor
s^{IJ}	2×2 matrix $s \equiv \text{diag}(1, -1)$
G	Coset space representative
L^μ	Cartan one-forms
L^α	Cartan superconnections
H_{LC}	The full lightcone string Hamiltonian
H_{int}	Interaction sector of the string Hamiltonian, appearing at $O(1/\widehat{R}^2)$ in the curvature expansion
H_{BB}	Purely bosonic sector of the interaction Hamiltonian

H_{FF}	Purely fermionic sector of the interaction Hamiltonian
H_{BF}	Bose-fermi mixing sector of the interaction Hamiltonian
Λ	String energy shift at $O(1/J)$; the shifts denoted by Λ_{BB} , Λ_{FF} and Λ_{BF} are associated with the corresponding sectors of H_{int}
ω_n	String energy at mode number n : $\sqrt{n^2 + p_-^2}$
k_n	String mode function $k_n = n$ (integer)
p_-	Worldsheet momentum in the x^- direction, equal to $1/\sqrt{\lambda'}$
L	Marks the overall level within a (super)multiplet of energy states
L_{sub}	Marks the level within an energy submultiplet (Chapter 5)
L'	Index labeling submultiplets within a supermultiplet of energy levels ($L = L_{\text{sub}} + L'$)

The various indices on the string side are chosen to represent the following:

$\mu, \nu, \rho = 0, \dots, 9$	$SO(9, 1)$ vectors ,
$\alpha, \beta, \gamma, \delta = 1, \dots, 16$	$SO(9, 1)$ spinors ,
$A, B = 1, \dots, 8$	$SO(8)$ vectors ,
$i, j, k = 1, \dots, 4$	$SO(4)$ vectors ,
$i', j', k' = 5, \dots, 8$	$SO(4)'$ vectors ,
$a, b = 0, 1$	Worldsheet coordinates (τ, σ) ,
$I, J, K, L = 1, 2$	Label two Majorana-Weyl spinors of equal chirality .

The 32×32 Dirac gamma matrices are decomposed into a 16×16 representation according to

$$\begin{aligned}
 (\Gamma^\mu)_{32 \times 32} &= \begin{pmatrix} 0 & \gamma^\mu \\ \bar{\gamma}^\mu & 0 \end{pmatrix} , & \gamma^\mu \bar{\gamma}^\nu + \gamma^\nu \bar{\gamma}^\mu &= 2\eta^{\mu\nu} , \\
 \gamma^\mu &= (1, \gamma^A, \gamma^9) , & \bar{\gamma}^\mu &= (-1, \gamma^A, \gamma^9) , \\
 \gamma^+ &= 1 + \gamma^9 , & \bar{\gamma}^+ &= -1 + \gamma^9 .
 \end{aligned} \tag{A.1}$$

In particular, the notation $\bar{\gamma}^\mu$ lowers the $SO(9, 1)$ spinor indices α, β :

$$\gamma^\mu = (\gamma^\mu)^{\alpha\beta} , \quad \bar{\gamma}^\mu = (\gamma^\mu)_{\alpha\beta} . \quad (\text{A.2})$$

These conventions are chosen to match those of Metsaev in [25]. By invoking κ -symmetry,

$$\bar{\gamma}^+ \theta = 0 \implies \bar{\gamma}^9 \theta = \theta , \quad (\text{A.3})$$

$$\bar{\gamma}^- = 1 + \bar{\gamma}^9 \implies \bar{\gamma}^- \theta = 2\theta . \quad (\text{A.4})$$

The antisymmetric product $\gamma^{\mu\nu}$ is given by

$$\begin{aligned} (\gamma^{\mu\nu})^\alpha_\beta &\equiv \frac{1}{2}(\gamma^\mu \bar{\gamma}^\nu)^\alpha_\beta - (\mu \rightleftharpoons \nu) , \\ (\bar{\gamma}^{\mu\nu})^\alpha_\beta &\equiv \frac{1}{2}(\bar{\gamma}^\mu \gamma^\nu)^\alpha_\beta - (\mu \rightleftharpoons \nu) . \end{aligned} \quad (\text{A.5})$$

We form the matrices Π and $\tilde{\Pi}$ according to:

$$\begin{aligned} \Pi &\equiv \gamma^1 \bar{\gamma}^2 \gamma^3 \bar{\gamma}^4 , \\ \tilde{\Pi} &\equiv \gamma^5 \bar{\gamma}^6 \gamma^7 \bar{\gamma}^8 . \end{aligned} \quad (\text{A.6})$$

These form the projection operators ($\Pi^2 = \tilde{\Pi}^2 = 1$)

$$\begin{aligned} \Pi_+ &\equiv \frac{1}{2}(1 + \Pi) , & \Pi_- &\equiv \frac{1}{2}(1 - \Pi) , \\ \tilde{\Pi}_+ &\equiv \frac{1}{2}(1 + \tilde{\Pi}) , & \tilde{\Pi}_- &\equiv \frac{1}{2}(1 - \tilde{\Pi}) . \end{aligned} \quad (\text{A.7})$$

The spinors θ^I represent two 32-component Majorana-Weyl spinors of $SO(9, 1)$ with equal chirality. The 32-component Weyl condition is $\Gamma_{11}\theta = \theta$, with

$$\Gamma_{11} = \Gamma^0 \dots \Gamma^9 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}_{32 \times 32} . \quad (\text{A.8})$$

The Weyl condition is used to select the top 16 components of θ to form the 16-component spinors

$$\theta^I = \begin{pmatrix} \theta^\alpha \\ 0 \end{pmatrix}^I . \quad (\text{A.9})$$

It is useful to form a single complex 16-component spinor ψ from the real spinors θ^1 and θ^2 :

$$\psi = \sqrt{2}(\theta^1 + i\theta^2) . \quad (\text{A.10})$$

The 16-component Weyl condition $\gamma^9\theta = \theta$ selects the upper eight components of θ , with

$$\gamma^9 = \gamma^1 \dots \gamma^8 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}_{16 \times 16} . \quad (\text{A.11})$$

The 16-component Dirac matrices γ^μ can, in turn, be constructed from the familiar Spin(8) Clifford algebra, wherein (in terms of $SO(8)$ vector indices)

$$(\gamma^A)_{16 \times 16} = \begin{pmatrix} 0 & \gamma^A \\ (\gamma^A)^T & 0 \end{pmatrix} , \quad (\text{A.12})$$

and

$$\{\gamma^A, \gamma^B\}_{16 \times 16} = 2\delta^{AB} , \quad (\gamma^A(\gamma^B)^T + \gamma^B(\gamma^A)^T = 2\delta^{AB})_{8 \times 8} . \quad (\text{A.13})$$

The Spin(8) Clifford algebra may be constructed explicitly in terms of eight real matrices

$$\begin{aligned} \gamma^1 &= \epsilon \times \epsilon \times \epsilon , & \gamma^5 &= \tau_3 \times \epsilon \times 1 , \\ \gamma^2 &= 1 \times \tau_1 \times \epsilon , & \gamma^6 &= \epsilon \times 1 \times \tau_1 , \end{aligned}$$

$$\begin{aligned}
\gamma^3 &= 1 \times \tau_3 \times \epsilon , & \gamma^7 &= \epsilon \times 1 \times \tau_3 , \\
\gamma^4 &= \tau_1 \times \epsilon \times 1 , & \gamma^8 &= 1 \times 1 \times 1 ,
\end{aligned} \tag{A.14}$$

with

$$\epsilon = \begin{pmatrix} 0 & 1 \\ -1 & 0 \end{pmatrix} , \quad \tau_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} , \quad \tau_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} . \tag{A.15}$$

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